LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100121\
Data File : BP007287.D
Acq On : 01 Oct 2021 12:46
Operator : CG/JU
Sample : PB139478BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : PB139478BS

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP007287.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.334	38	42	53	rVB	321137	417353	3.05%	0.125%
2	3.652	92	96	109	rBV	508104	811774	5.93%	0.244%
3	3.746	109	112	146	rVB	549579	1066404	7.79%	0.320%
4	4.981	317	322	334	rBV	661253	945319	6.90%	0.284%
5	5.452	395	402	420	rBV	3381573	5004552	36.55%	1.502%
6	7.010	659	667	669	rBV	780354	1182017	8.63%	0.355%
7	7.052	669	674	691	rVV2	3055366	6932719	50.63%	2.081%
8	7.199	691	699	709	rVV	1091163	1821799	13.31%	0.547%
9	7.293	709	715	724	rVB	974272	1417605	10.35%	0.425%
10	7.434	731	739	751	rBV	1283581	1985328	14.50%	0.596%
11	7.752	786	793	805	rBV	1429937	2244049	16.39%	0.673%
12	7.863	805	812	815	rVV	770223	1287320	9.40%	0.386%
13	7.899	815	818	830	rVB	1462872	2232202	16.30%	0.670%
14	8.116	845	855	865	rBV	1371001	2294498	16.76%	0.689%
15	8.216	865	872	883	rVB	1406562	2279630	16.65%	0.684%
16	8.322	883	890	896	rBV	1526666	2391698	17.47%	0.718%
17	8.393	896	902	910	rVB2	654730	1597598	11.67%	0.479%
18	8.652	938	946	949	rBV	1370754	2677088	19.55%	0.803%
19	8.693	949	953	973	rVB2	1953934	3814669	27.86%	1.145%
20	8.940	988	995	1003	rVV	1165336	1904846	13.91%	0.572%
21	9.028	1003	1010	1014	rVV	1835867	3190523	23.30%	0.958%
22	9.075	1014	1018	1032	rVB	1059693	1716621	12.54%	0.515%
23	9.293	1049	1055	1066	rBV	91359	188980	1.38%	0.057%
24	9.604	1098	1108	1119	rBV	1383666	2283234	16.68%	0.685%
25	9.781	1130	1138	1144	rBV	877054	1453444	10.61%	0.436%
26	9.851	1144	1150	1157	rVB	1692556	2660895	19.43%	0.799%
27	10.028	1163	1180	1184	rBV	518625	1743310	12.73%	0.523%
28	10.081	1184	1189	1201	rVB	1170424	1833765	13.39%	0.550%
29	10.322	1221	1230	1248	rBV	1416757	2421475	17.68%	0.727%
30	10.528	1258	1265	1276	rVB	1628583	2728190	19.92%	0.819%
31	10.663	1279	1288	1292	rBV	1004565	1750980	12.79%	0.526%
32	10.716	1292	1297	1305	rVB	2071564	3375410	24.65%	1.013%
33	10.834	1308	1317	1332	rBV	1087372	1946480	14.22%	0.584%
34	10.998	1338	1345	1353	rVB	1565391	2585398	18.88%	0.776%
35	11.646	1446	1455	1459	rBV	355903	777043	5.68%	0.233%
36	11.969	1500	1510	1523	rBV	1610865	2528321	18.47%	0.759%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100121\
Data File : BP007287.D
Acq On : 01 Oct 2021 12:46
Operator : CG/JU
Sample : PB139478BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : PB139478BS

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	12.222	1547	1553	1563	rVB	337991	518184	3.78%	0.156%
38	12.328	1563	1571	1585	rBV	2573072	4030381	29.44%	1.210%
39	12.457	1586	1593	1600	rBV	106599	265227	1.94%	0.080%
40	12.545	1601	1608	1623	rVB	2498635	3835229	28.01%	1.151%
41	12.669	1623	1629	1632	rBV	2300481	4055938	29.62%	1.217%
42	12.940	1668	1675	1680	rBV	1800014	2725000	19.90%	0.818%
43	13.016	1680	1688	1699	rVV2	1661401	3463078	25.29%	1.039%
44	13.128	1699	1707	1720	rVB	5134318	7628001	55.71%	2.289%
45	13.334	1735	1742	1746	rBV	2834666	4243195	30.99%	1.273%
46	13.375	1746	1749	1758	rVB	2289141	3473797	25.37%	1.043%
47	13.587	1773	1785	1794	rBV	1248208	1976036	14.43%	0.593%
48	13.881	1829	1835	1843	rBV	985236	1397805	10.21%	0.420%
49	13.963	1843	1849	1855	rVV	2291153	3162316	23.10%	0.949%
50	14.028	1855	1860	1865	rVV	1078658	1606618	11.73%	0.482%
51	14.087	1865	1870	1878	rVV	1450281	2109306	15.40%	0.633%
52	14.187	1879	1887	1889	rVV	900790	1312436	9.59%	0.394%
53	14.222	1889	1893	1904	rVB	2644048	3874539	28.30%	1.163%
54	14.416	1920	1926	1933	rBV	1103899	1638856	11.97%	0.492%
55	14.498	1934	1940	1946	rVV	1563790	2298364	16.79%	0.690%
56	14.563	1946	1951	1957	rVV	3111077	4508204	32.92%	1.353%
57	14.628	1957	1962	1974	rVV	1645770	2451742	17.91%	0.736%
58	14.739	1974	1981	1987	rVV	1898187	3104955	22.68%	0.932%
59	14.798	1987	1991	1998	rVV	429587	706749	5.16%	0.212%
60	14.875	1998	2004	2005	rVV	1641650	2056788	15.02%	0.617%
61	14.898	2005	2008	2020	rVB	2833438	4034220	29.46%	1.211%
62	15.051	2027	2034	2042	rBV	1813370	2644426	19.31%	0.794%
63	15.134	2042	2048	2068	rVB	1927792	2839463	20.74%	0.852%
64	15.328	2074	2081	2086	rBV	2581485	3496640	25.54%	1.049%
65	15.545	2111	2118	2122	rBV2	5279659	8173371	59.69%	2.453%
66	15.581	2122	2124	2129	rVV	1217574	1513446	11.05%	0.454%
67	15.639	2129	2134	2144	rVV	1117642	1547993	11.31%	0.465%
68	15.763	2148	2155	2160	rVV	3346758	4500617	32.87%	1.351%
69	15.839	2162	2168	2176	rVB	2271836	3229414	23.59%	0.969%
70	15.992	2187	2194	2208	rBV	3909510	5751525	42.01%	1.726%
71	16.439	2263	2270	2277	rBV	2347560	3104618	22.67%	0.932%
72	16.557	2283	2290	2302	rVB	2533147	3502753	25.58%	1.051%
73	16.722	2305	2318	2332	rBV	2521995	3259208	23.80%	0.978%
74	16.910	2343	2350	2365	rBV	3304638	4965750	36.27%	1.490%
75	17.251	2402	2408	2411	rBV	1895142	2657282	19.41%	0.798%
76	17.298	2411	2416	2422	rVV	3745039	5382607	39.31%	1.615%

LSC Area Percent Report

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Operator : CG/JU
Sample : PB139478BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : PB139478BS

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	17.386	2422	2431	2443	rVB	3803855	5321079	38.86%	1.597%
78	17.663	2472	2478	2491	rVB	3365818	4714918	34.43%	1.415%
79	18.210	2565	2571	2576	rBV	3737607	4345463	31.74%	1.304%
80	19.263	2744	2750	2758	rBV	4894955	6094186	44.51%	1.829%
81	19.445	2775	2781	2787	rBV	2415682	3130005	22.86%	0.939%
82	19.510	2787	2792	2797	rVB2	111481	208601	1.52%	0.063%
83	19.616	2802	2810	2820	rVV	4808858	6388567	46.66%	1.917%
84	19.810	2837	2843	2848	rBV	9370933	11300791	82.53%	3.392%
85	20.051	2880	2884	2891	rBV	561652	722375	5.28%	0.217%
86	20.486	2953	2958	2962	rBV	3971319	4470656	32.65%	1.342%
87	20.551	2962	2969	2974	rVV	4605657	5310806	38.79%	1.594%
88	20.704	2991	2995	2999	rBV	134287	172357	1.26%	0.052%
89	21.251	3083	3088	3095	rBV2	5857475	9102256	66.48%	2.732%
90	21.327	3095	3101	3106	rVV2	5498415	9681584	70.71%	2.906%
91	21.374	3106	3109	3116	rVB	5315704	6641531	48.51%	1.993%
92	21.610	3146	3149	3156	rVB3	167713	210960	1.54%	0.063%
93	22.168	3239	3244	3250	rVB	4087287	5437834	39.71%	1.632%
94	23.010	3380	3387	3391	rBV	3903360	7213282	52.68%	2.165%
95	23.057	3391	3395	3405	rVB	3892534	6629800	48.42%	1.990%
96	23.639	3486	3494	3503	rVB	3669516	7417615	54.17%	2.226%
97	23.745	3505	3512	3525	rVB2	1421263	3029866	22.13%	0.909%
98	24.357	3612	3616	3624	rVB	401735	799843	5.84%	0.240%
99	26.274	3929	3942	3955	rVB3	3903821	13692378	100.00%	4.109%
100	27.039	4062	4072	4088	rVB	2003964	6617492	48.33%	1.986%

Sum of corrected areas: 333192859

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100121\
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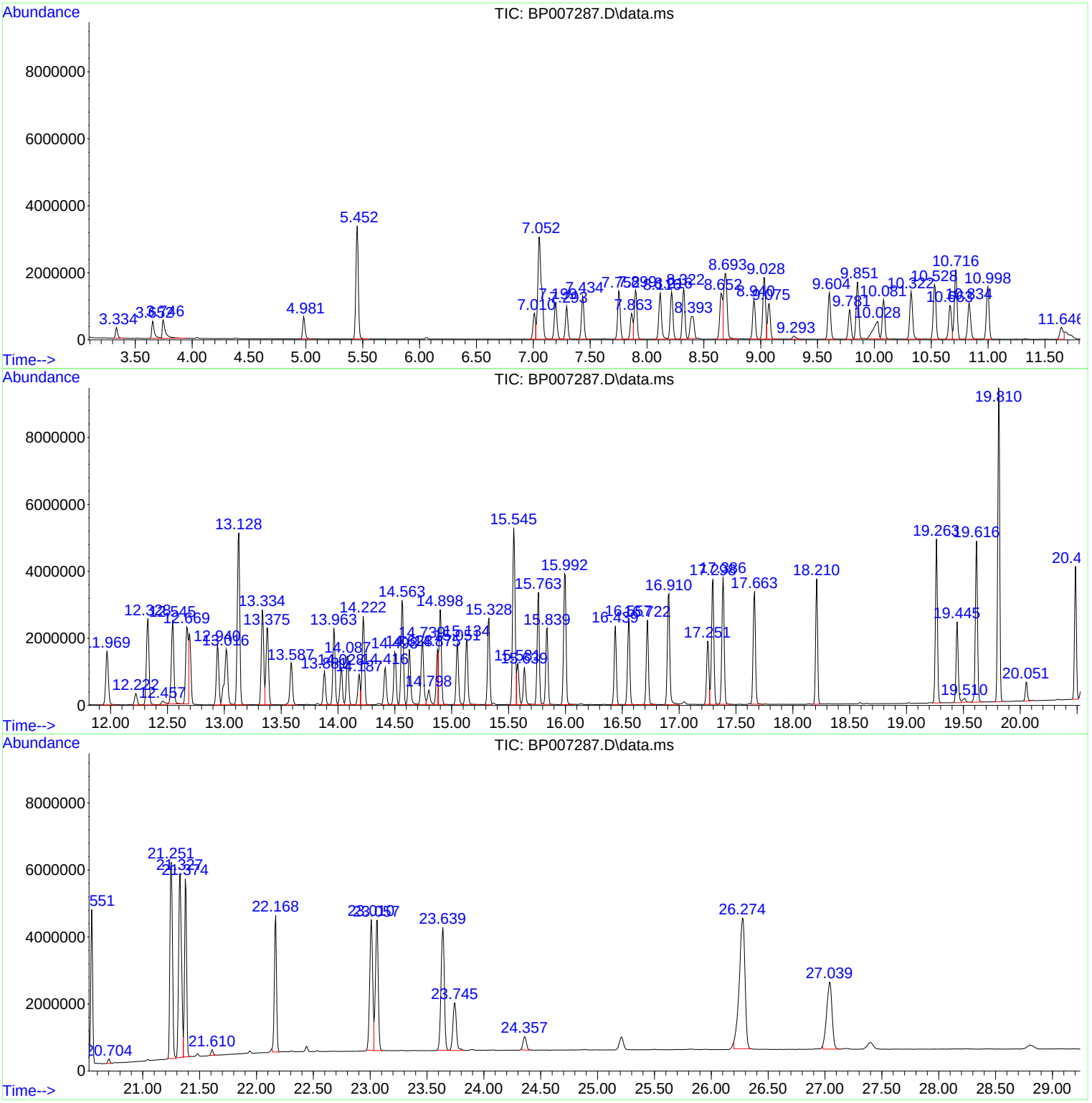
Sample : PB139478BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB139478BS

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100121\
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Instrument : BNA_P
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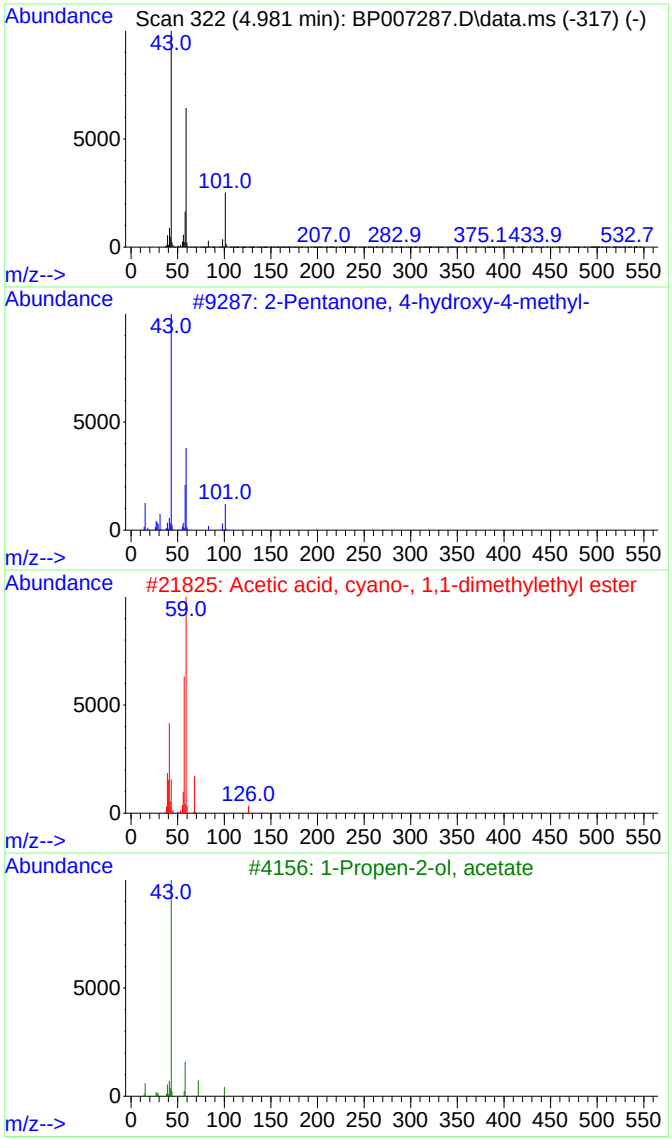
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

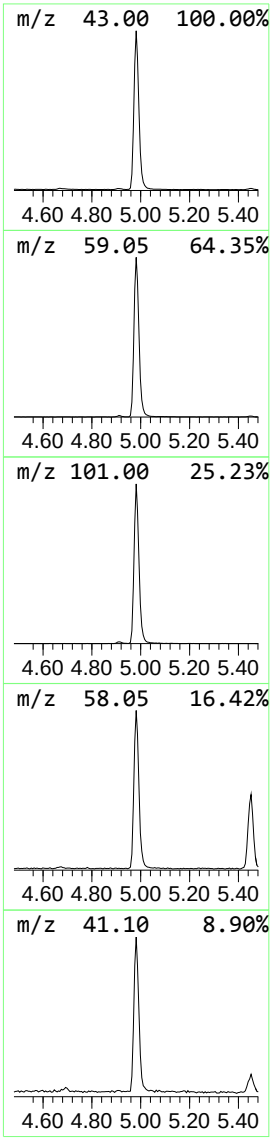
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.981	14.69 ng	945319	1,4-Dichlorobenzene-d4	7.863

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100121\
Data File : BP007287.D
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ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB139478BS

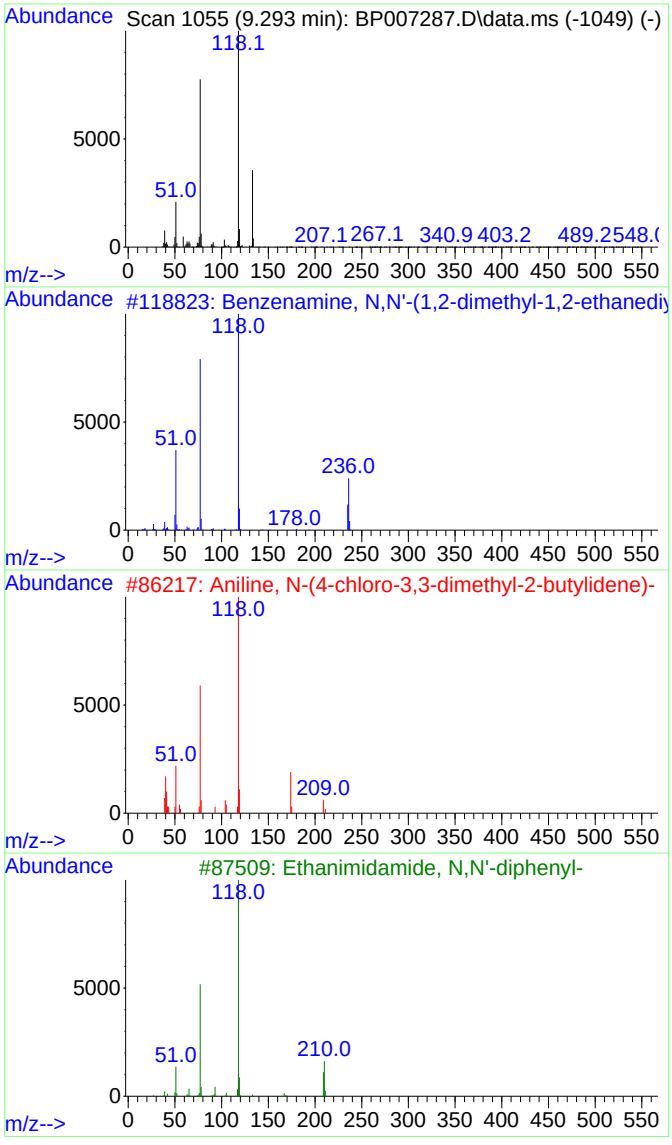
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

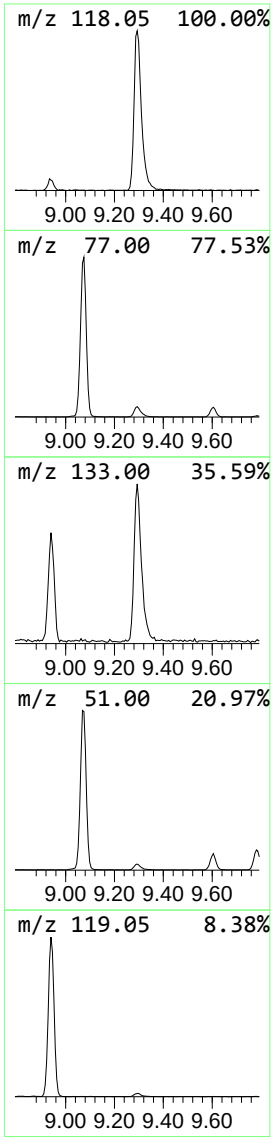
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzenamine, N,N'-(1,2-dime... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.293	2.16 ng	188980	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, N,N'-(1,2-dimethyl-...	236	C16H16N2	005393-49-7	64
2	Aniline, N-(4-chloro-3,3-dimethy...	209	C12H16ClN	1010137-59-6	64
3	Ethanimidamide, N,N'-diphenyl-	210	C14H14N2	000621-09-0	39
4	Methanamine, N-(1-phenylethylide...	133	C9H11N	006907-71-7	38
5	(1-Phenylethylidene)[(trimethyls...	207	C11H17NOSi	064948-39-6	33





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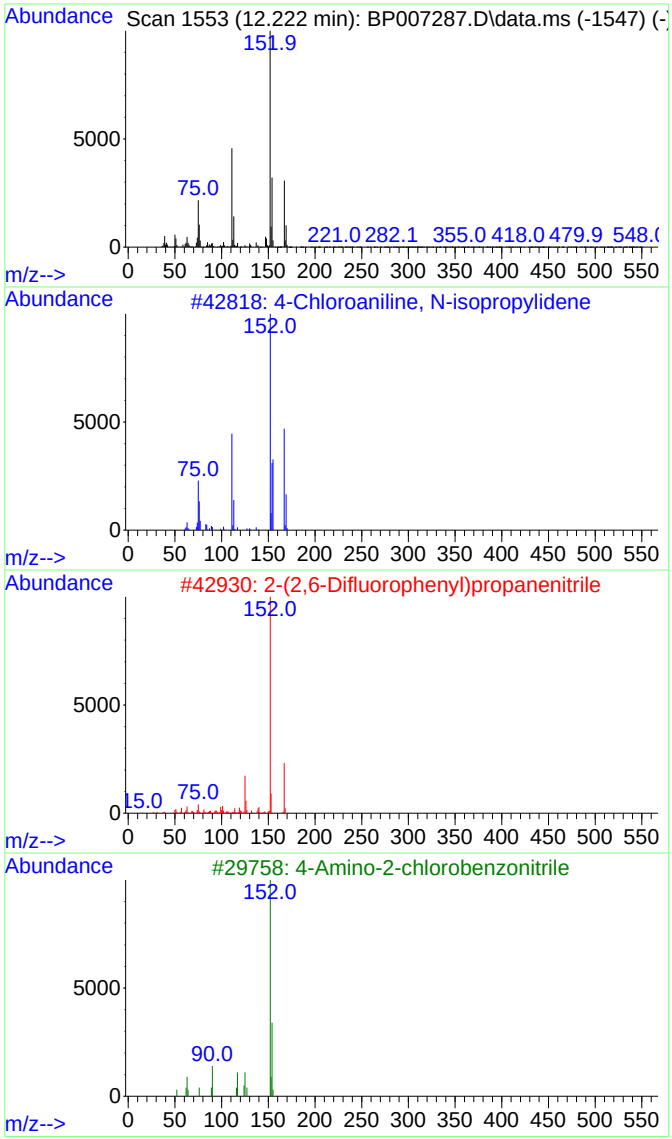
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

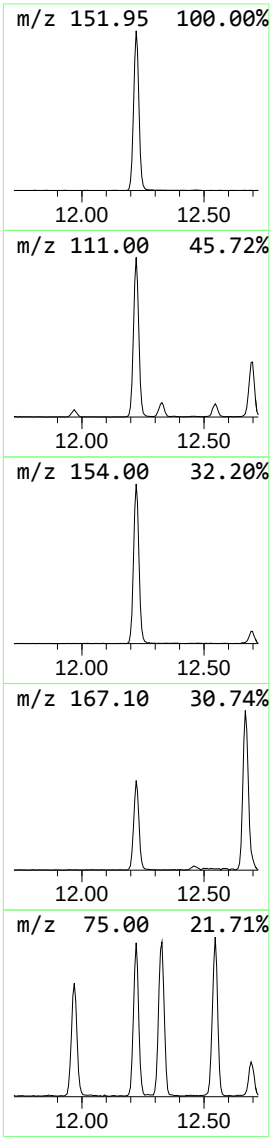
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 4-Chloroaniline, N-isopropy... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.222	5.92 ng	518184	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Chloroaniline, N-isopropylidene	167	C9H10ClN	040938-43-0	95
2	2-(2,6-Difluorophenyl)propanenit...	167	C9H7F2N	149680-18-2	38
3	4-Amino-2-chlorobenzonitrile	152	C7H5ClN2	020925-27-3	38
4	5-Aminopyrimidine, N-trimethylsi...	167	C7H13N3Si	1000493-88-4	38
5	Chlorhexidine	504	C22H30Cl2N10	000055-56-1	38





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Sample : PB139478BS
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Instrument :
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ClientSampleId :
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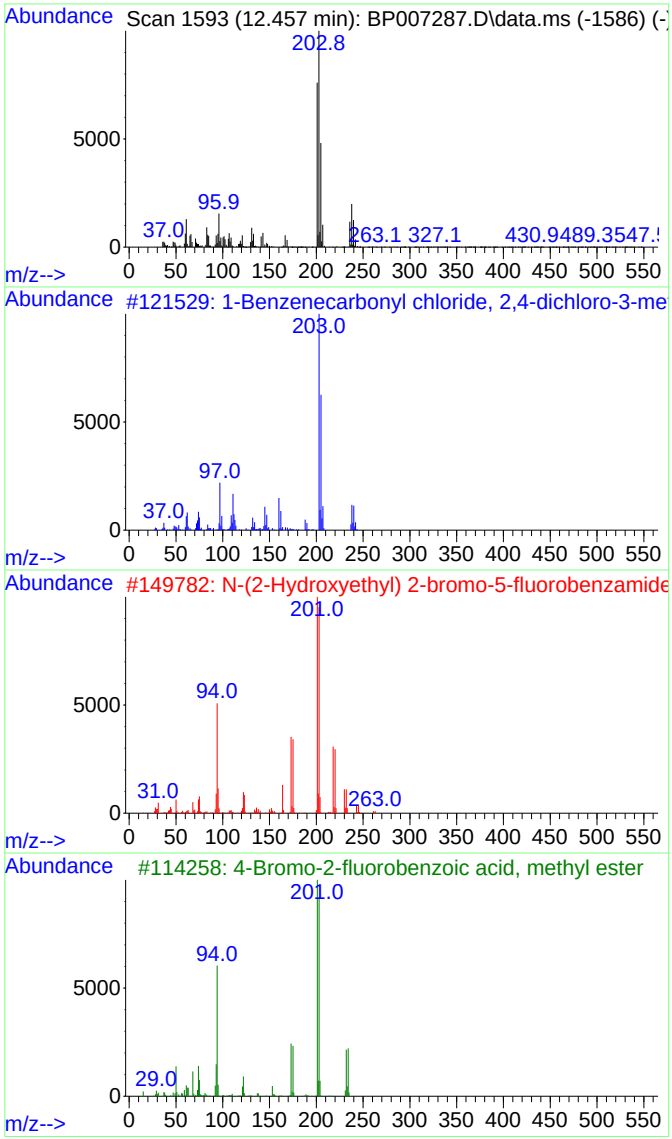
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

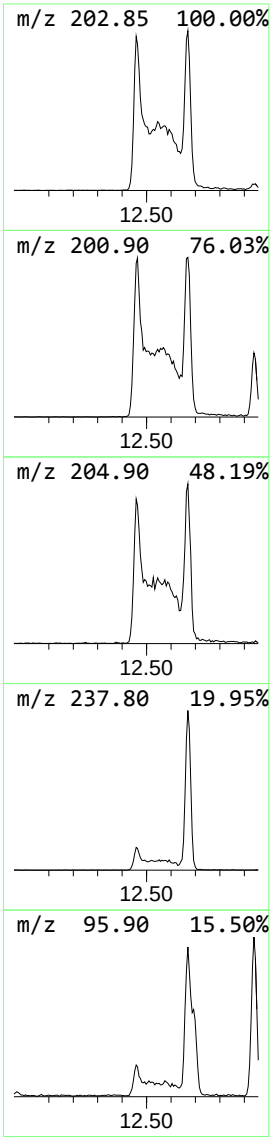
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 1-Benzenecarbonyl chloride,... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.457	3.03 ng	265227	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Benzenecarbonyl chloride, 2,4-...	238	C8H5Cl3O2	1000196-85-1	46
2	N-(2-Hydroxyethyl) 2-bromo-5-flu...	261	C9H9BrFNO2	951884-16-5	38
3	4-Bromo-2-fluorobenzoic acid, me...	232	C8H6BrFO2	179232-29-2	37
4	5-Bromonicotinic acid	201	C6H4BrNO2	020826-04-4	37
5	Pyridine, 2,4-dichloro-3,5,6-tri...	201	C5Cl2F3N	054889-43-9	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100121\
Data File : BP007287.D
Acq On : 01 Oct 2021 12:46
Operator : CG/JU
Sample : PB139478BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB139478BS

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

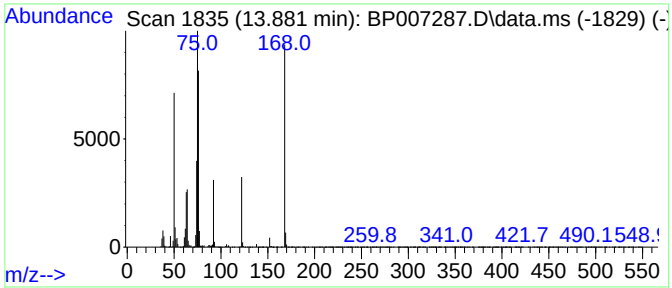
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

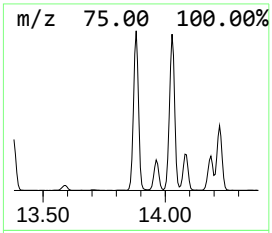
****

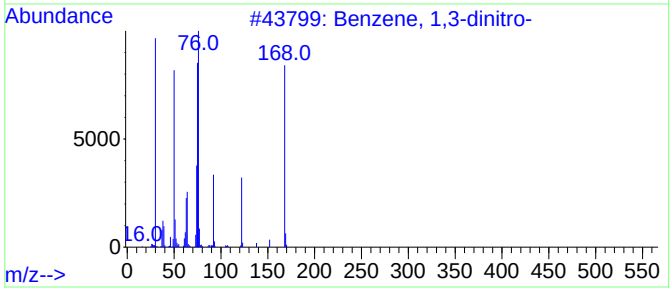
Peak Number 5 Benzene, 1,3-dinitro- Concentration Rank 4

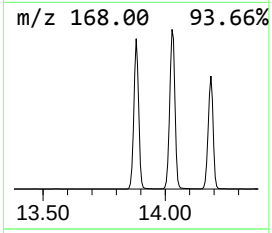
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.881	12.16 ng	1397810	Acenaphthene-d10	14.498

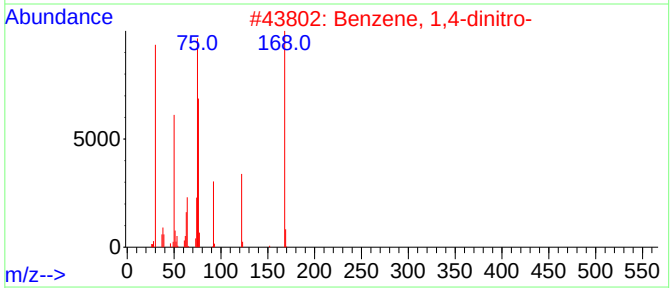
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dinitro-	168	C6H4N2O4	000099-65-0	98
2	Benzene, 1,4-dinitro-	168	C6H4N2O4	000100-25-4	97
3	Bicyclo[4.2.0]octa-1,3,5-triene-...	132	C8H4O2	006383-11-5	33
4	1H-Isoindole-1,3(2H)-dione, 2-(h...	177	C9H7NO3	000118-29-6	17
5	Benzaldehyde, 2,6-dinitro-	196	C7H4N2O5	000606-31-5	17

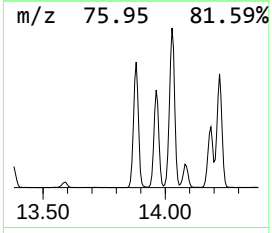


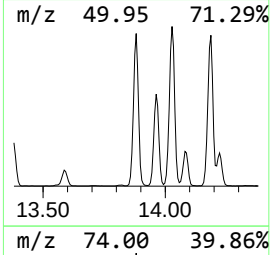


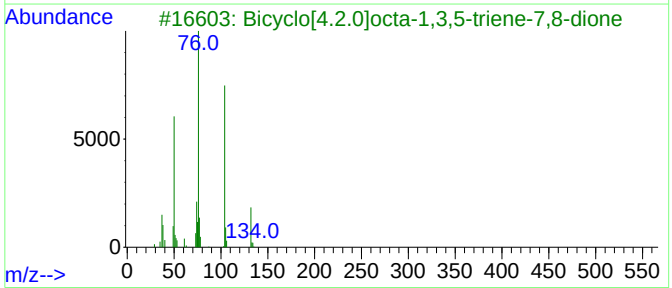


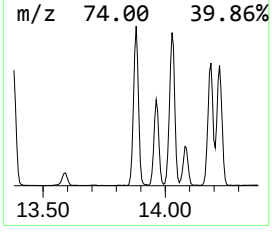












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100121\
Data File : BP007287.D
Acq On : 01 Oct 2021 12:46
Operator : CG/JU
Sample : PB139478BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB139478BS

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

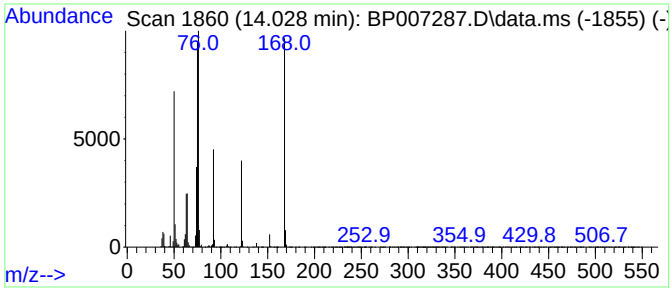
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

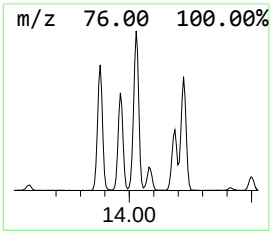
****

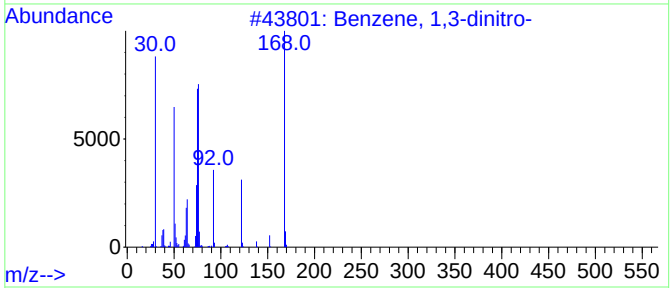
Peak Number 6 Benzene, 1,3-dinitro- Concentration Rank 3

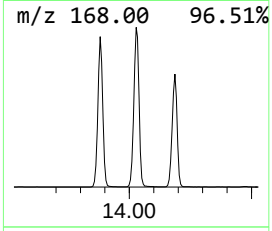
R.T.	EstConc	Area	Relative to ISTD	R.T.
14.028	13.98 ng	1606620	Acenaphthene-d10	14.498

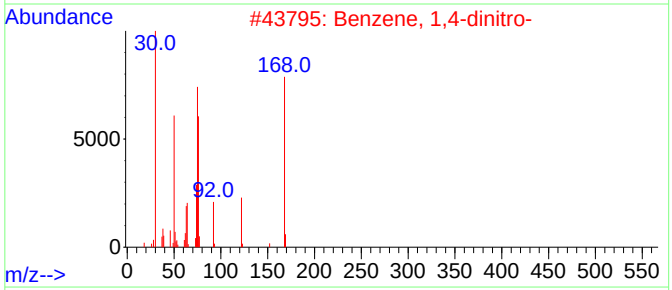
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dinitro-	168	C6H4N2O4	000099-65-0	98
2	Benzene, 1,4-dinitro-	168	C6H4N2O4	000100-25-4	94
3	1H-Isoindole-1,3(2H)-dione, 2-(h...	177	C9H7NO3	000118-29-6	17
4	Methyl hydrogen phthalate	180	C9H8O4	004376-18-5	9
5	Bicyclo[4.2.0]octa-1,3,5-triene-...	132	C8H4O2	006383-11-5	9

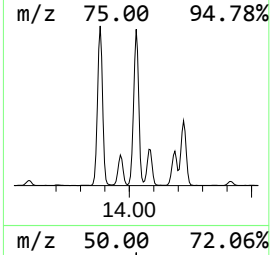


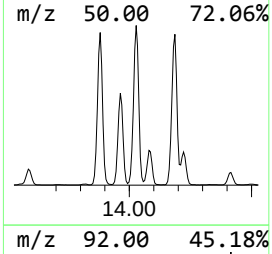


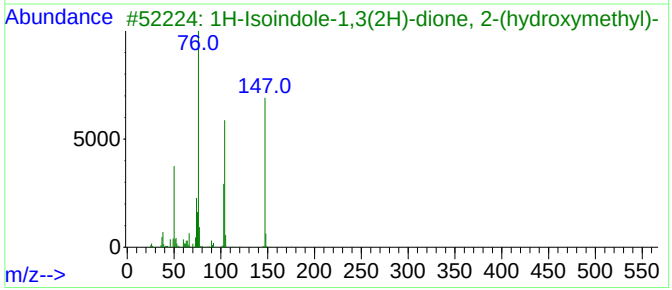


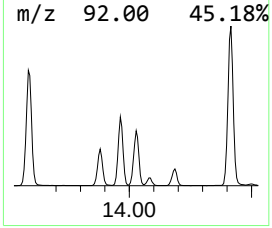












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100121\
Data File : BP007287.D
Acq On : 01 Oct 2021 12:46
Operator : CG/JU
Sample : PB139478BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : PB139478BS

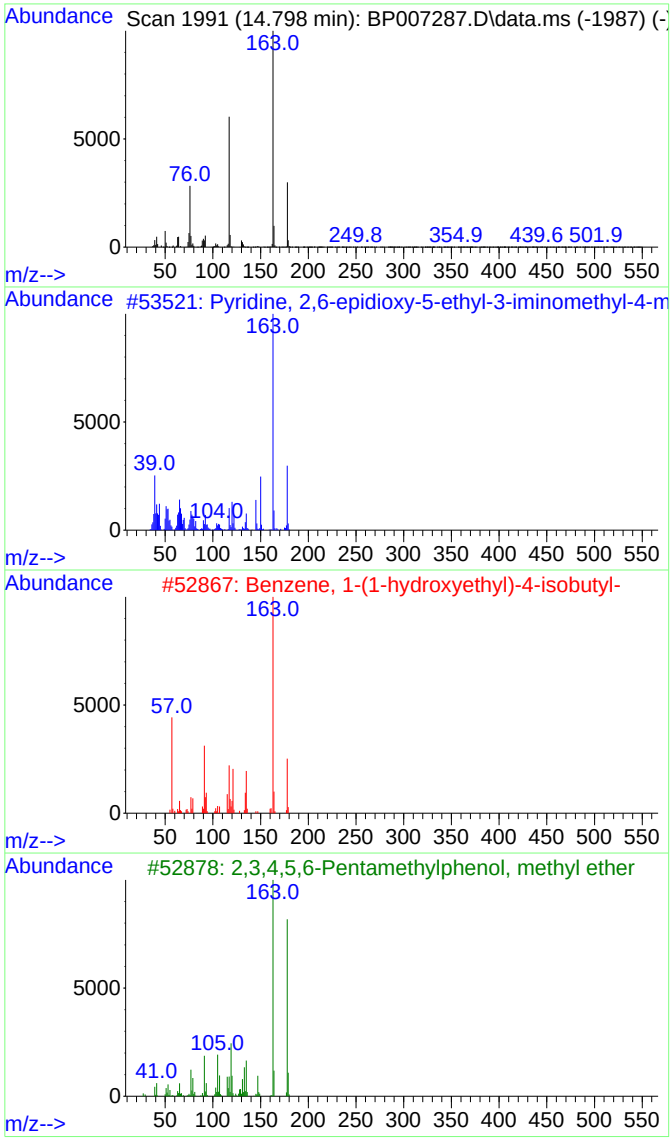
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

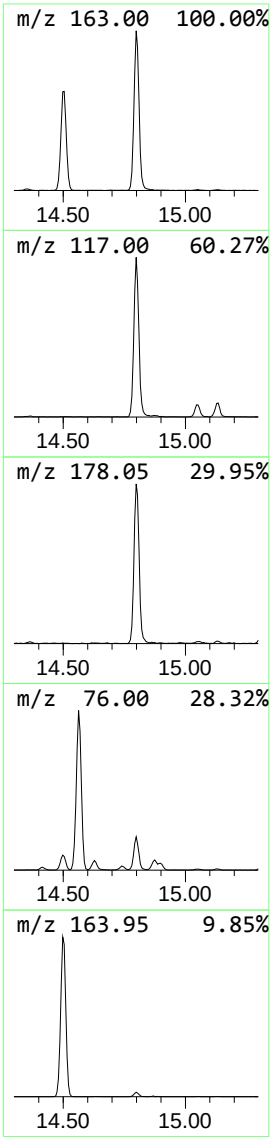
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Pyridine, 2,6-epidioxy-5-et... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.798	6.15 ng	706749	Acenaphthene-d10	14.498

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyridine, 2,6-epidioxy-5-ethyl-3...	178	C9H10N2O2	1000263-31-1	50
2	Benzene, 1-(1-hydroxyethyl)-4-is...	178	C12H18O	040150-92-3	50
3	2,3,4,5,6-Pentamethylphenol, met...	178	C12H18O	014804-37-6	47
4	Propofol	178	C12H18O	002078-54-8	47
5	Ethanone, 1-(2,3-dihydro-1,4-ben...	178	C10H10O3	002879-20-1	46





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100121\
Data File : BP007287.D
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Operator : CG/JU
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Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB139478BS

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

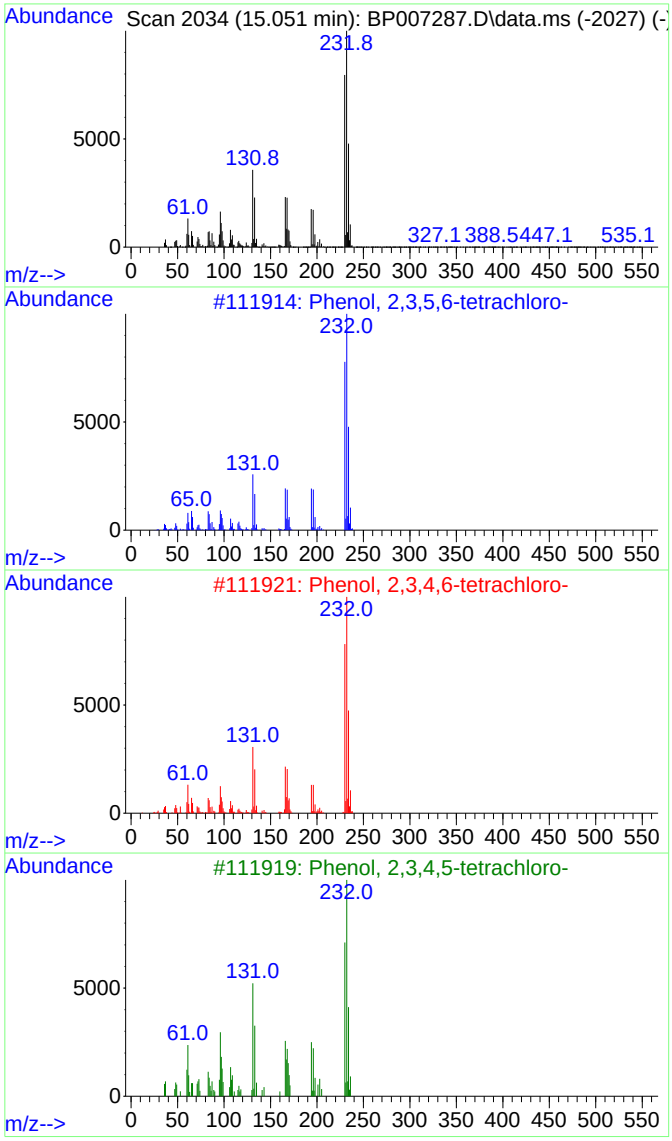
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

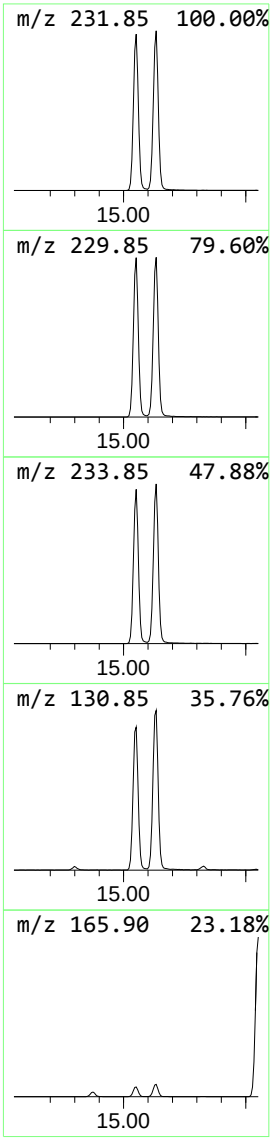
****

Peak Number 8 Phenol, 2,3,5,6-tetrachloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.051	23.01 ng	2644430	Acenaphthene-d10	14.498

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	99
2	Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	99
3	Phenol, 2,3,4,5-tetrachloro-	230	C6H2Cl4O	004901-51-3	93
4	4-Aminotetrachloropyridine	230	C5H2Cl4N2	002176-63-8	68
5	2,3,4,6-Tetrachlorophenol, isopr...	272	C9H8Cl4O	1010395-04-9	64





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Data File : BP007287.D
Acq On : 01 Oct 2021 12:46
Operator : CG/JU
Sample : PB139478BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : PB139478BS

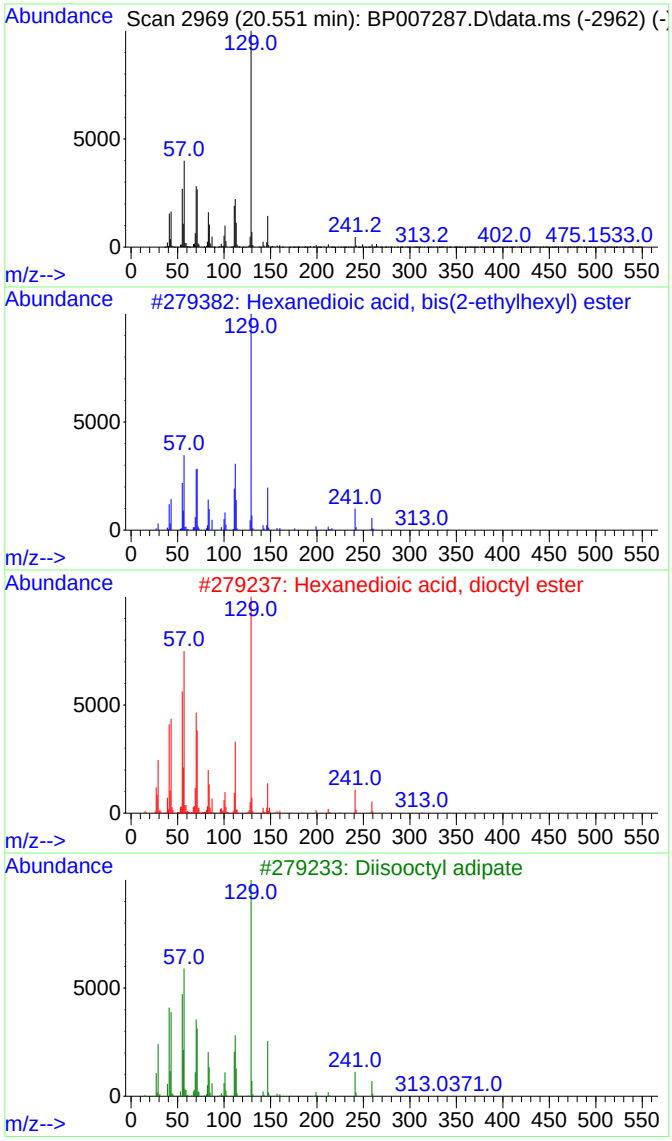
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

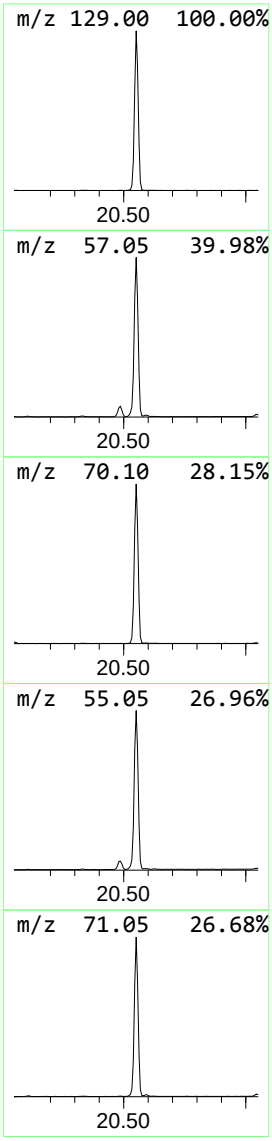
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Hexanedioic acid, bis(2-eth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.551	10.97 ng	5310810	Chrysene-d12	21.339

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	91
2	Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	87
3	Diisooctyl adipate	370	C22H42O4	001330-86-5	78
4	Hexanedioic acid, mono(2-ethylhe...	258	C14H26O4	004337-65-9	68
5	Adipic acid, 2-ethylhexyl isobut...	314	C18H34O4	1010324-48-0	64





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Data File : BP007287.D
Acq On : 01 Oct 2021 12:46
Operator : CG/JU
Sample : PB139478BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : PB139478BS

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

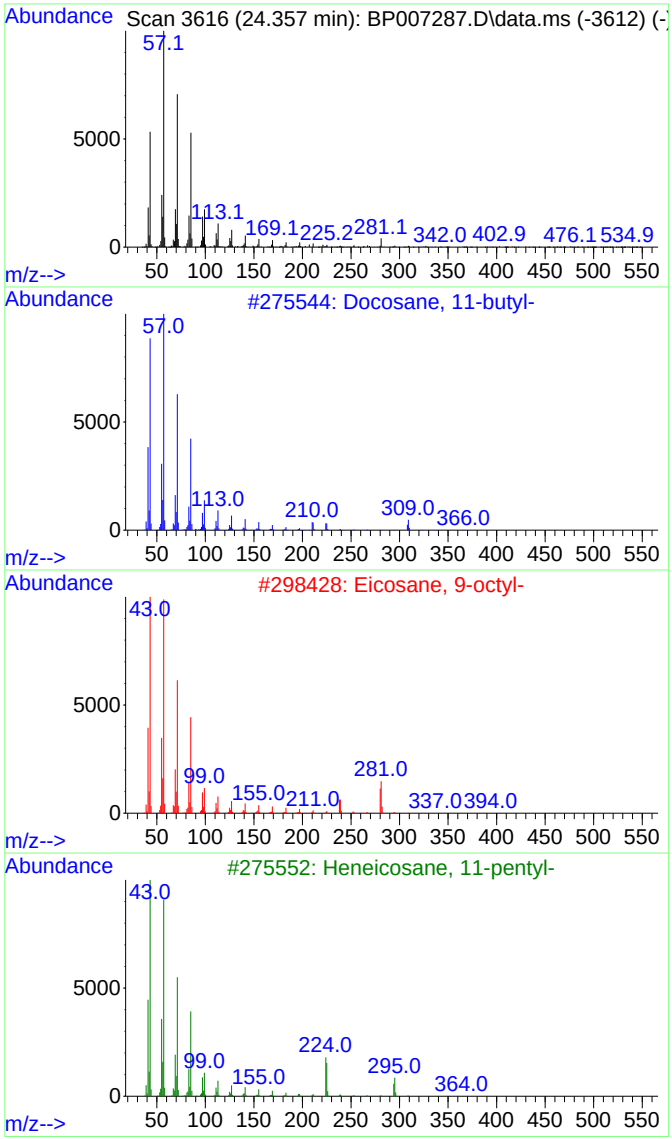
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

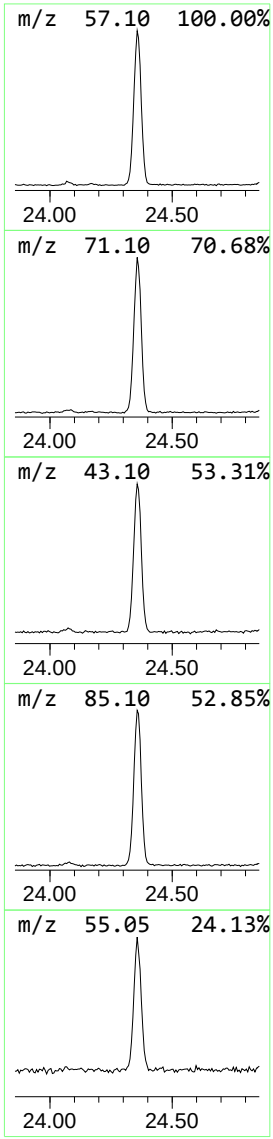
****

Peak Number 10 Docosane, 11-butyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.357	5.28 ng	799843	Perylene-d12	23.745

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-	366	C26H54	013475-76-8	93
2	Eicosane, 9-octyl-	394	C28H58	013475-77-9	93
3	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	93
4	Docosane, 5-butyl-	366	C26H54	055282-16-1	91
5	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100121\
Data File : BP007287.D
Acq On : 01 Oct 2021 12:46
Operator : CG/JU
Sample : PB139478BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : PB139478BS

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Pentanone, 4-...	4.981	14.7	ng	945319	1	7.863	1287320	20.0
Benzenamine, N,...	9.293	2.2	ng	188980	2	10.663	1750980	20.0
4-Chloroaniline...	12.222	5.9	ng	518184	2	10.663	1750980	20.0
1-Benzenecarbon...	12.457	3.0	ng	265227	2	10.663	1750980	20.0
Benzene, 1,3-di...	13.881	12.2	ng	1397810	3	14.498	2298360	20.0
Benzene, 1,3-di...	14.028	14.0	ng	1606620	3	14.498	2298360	20.0
Pyridine, 2,6-e...	14.798	6.2	ng	706749	3	14.498	2298360	20.0
Phenol, 2,3,5,6...	15.051	23.0	ng	2644430	3	14.498	2298360	20.0
Hexanedioic aci...	20.551	11.0	ng	5310810	5	21.339	9681580	20.0
Docosane, 11-bu...	24.357	5.3	ng	799843	6	23.745	3029870	20.0