LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010914.D
Acq On : 30 Jun 2022 11:30
Operator : CG/JU
Sample : N3523-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JT6

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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\SFAM-EPA-BP062322.M
Title : SVOA CALIBRATION

Signal : TIC: BP010914.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.205	15	20	28	rVB	133835	175257	0.98%	0.121%
2	3.476	62	66	71	rBV	57046	72844	0.41%	0.050%
3	3.611	87	89	106	rBV	535897	787861	4.42%	0.545%
4	3.834	119	127	133	rBV	47505	76353	0.43%	0.053%
5	3.928	138	143	148	rBV	46410	62946	0.35%	0.044%
6	5.034	323	331	343	rBV	1986820	2848305	15.99%	1.969%
7	5.170	349	354	363	rBV5	8694	18109	0.10%	0.013%
8	5.428	394	398	404	rVB	13520	21014	0.12%	0.015%
9	5.770	452	456	459	rBV6	11937	20009	0.11%	0.014%
10	5.805	459	462	471	rVB5	17136	27810	0.16%	0.019%
11	6.881	634	645	658	rVB	461882	795660	4.47%	0.550%
12	7.046	667	673	681	rBV	1849475	2794823	15.69%	1.932%
13	7.122	681	686	694	rVB	32512	54279	0.30%	0.038%
14	7.240	698	706	717	rBV	1727952	2773428	15.57%	1.917%
15	7.328	717	721	728	rVV6	12520	18110	0.10%	0.013%
16	7.428	732	738	744	rVB	34731	58018	0.33%	0.040%
17	7.593	758	766	776	rVB	553274	840577	4.72%	0.581%
18	7.705	776	785	788	rBV	1859992	2987167	16.77%	2.065%
19	7.740	788	791	806	rVB	1649717	2392031	13.43%	1.653%
20	8.058	836	845	853	rBV	11143454	17811640	100.00%	12.311%
21	8.422	899	907	918	rVV	1285031	2012120	11.30%	1.391%
22	8.534	920	926	936	rVB5	15526	33192	0.19%	0.023%
23	8.869	975	983	986	rBV	1896282	3198644	17.96%	2.211%
24	8.911	986	990	1000	rVV	3735368	5728806	32.16%	3.960%
25	8.993	1000	1004	1008	rVV3	18116	31791	0.18%	0.022%
26	9.199	1034	1039	1044	rBV5	11405	18093	0.10%	0.013%
27	9.287	1049	1054	1058	rBV	104935	167523	0.94%	0.116%
28	9.511	1085	1092	1097	rBV2	35661	60452	0.34%	0.042%
29	9.587	1097	1105	1119	rVV	1360704	2352960	13.21%	1.626%
30	10.122	1189	1196	1204	rBV	2022497	3325660	18.67%	2.299%
31	10.187	1204	1207	1214	rVV2	43156	69490	0.39%	0.048%
32	10.281	1215	1223	1229	rVB3	39877	80788	0.45%	0.056%
33	10.363	1229	1237	1252	rBV	4726501	7517937	42.21%	5.196%
34	10.499	1252	1260	1267	rBV	2502748	4030247	22.63%	2.786%
35	10.646	1275	1285	1296	rBV	1439810	2461010	13.82%	1.701%
36	10.881	1317	1325	1336	rVB	2304559	3785806	21.25%	2.617%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010914.D
Acq On : 30 Jun 2022 11:30
Operator : CG/JU
Sample : N3523-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JT6

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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\SFAM-EPA-BP062322.M
Title : SVOA CALIBRATION

37	11.105	1352	1363	1371	rBV	2374325	3757323	21.09%	2.597%
38	11.322	1392	1400	1410	rVB	534034	897265	5.04%	0.620%
39	11.605	1442	1448	1453	rBV2	16505	27411	0.15%	0.019%
40	11.799	1471	1481	1490	rBV2	238700	442597	2.48%	0.306%
41	12.093	1526	1531	1537	rVB	44674	69841	0.39%	0.048%
42	12.487	1590	1598	1603	rBV	195946	329928	1.85%	0.228%
43	12.534	1603	1606	1617	rVB	145298	229438	1.29%	0.159%
44	12.781	1641	1648	1656	rVV3	94714	172673	0.97%	0.119%
45	12.869	1656	1663	1669	rVV10	12216	27278	0.15%	0.019%
46	12.934	1669	1674	1678	rVV	75347	111318	0.62%	0.077%
47	12.975	1678	1681	1690	rVB8	21152	49275	0.28%	0.034%
48	13.152	1700	1711	1724	rBV	835035	1422408	7.99%	0.983%
49	13.263	1724	1730	1735	rBV4	33200	51235	0.29%	0.035%
50	13.322	1735	1740	1744	rVB3	39646	60995	0.34%	0.042%
51	13.381	1744	1750	1757	rBV2	152842	243327	1.37%	0.168%
52	13.781	1806	1818	1832	rBV	3863239	5492091	30.83%	3.796%
53	14.051	1851	1864	1874	rBV	4302966	6171174	34.65%	4.265%
54	14.363	1907	1917	1929	rBV	3131966	4704232	26.41%	3.251%
55	14.575	1948	1953	1962	rBV	259739	400571	2.25%	0.277%
56	14.675	1964	1970	1976	rVB4	32018	61144	0.34%	0.042%
57	14.851	1994	2000	2006	rBV	103772	160082	0.90%	0.111%
58	14.993	2020	2024	2028	rVB3	30025	44393	0.25%	0.031%
59	15.122	2042	2046	2051	rVB	24558	32690	0.18%	0.023%
60	15.181	2051	2056	2062	rBV	48749	67995	0.38%	0.047%
61	15.357	2080	2086	2092	rBV	5114705	7450251	41.83%	5.149%
62	15.404	2092	2094	2101	rVV	177534	241477	1.36%	0.167%
63	15.481	2101	2107	2121	rVV	1131815	1560726	8.76%	1.079%
64	15.598	2123	2127	2133	rVB	59795	89372	0.50%	0.062%
65	15.945	2180	2186	2191	rBV10	21620	44070	0.25%	0.030%
66	16.151	2217	2221	2228	rVB6	39937	63886	0.36%	0.044%
67	16.345	2248	2254	2260	rBV7	51153	99420	0.56%	0.069%
68	16.522	2280	2284	2290	rVB8	22195	32356	0.18%	0.022%
69	16.687	2308	2312	2318	rBV8	21610	37059	0.21%	0.026%
70	16.934	2352	2354	2363	rVB2	25967	36255	0.20%	0.025%
71	17.122	2380	2386	2397	rBV	3109403	4501997	25.28%	3.112%
72	17.222	2397	2403	2415	rVV2	5153306	7548249	42.38%	5.217%
73	17.328	2417	2421	2424	rVV5	38013	69093	0.39%	0.048%
74	17.363	2424	2427	2436	rVB4	41774	77265	0.43%	0.053%
75	17.492	2447	2449	2453	rBV5	11983	18451	0.10%	0.013%
76	17.534	2453	2456	2462	rVB7	17869	25701	0.14%	0.018%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010914.D
Acq On : 30 Jun 2022 11:30
Operator : CG/JU
Sample : N3523-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JT6

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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\SFAM-EPA-BP062322.M
Title : SVOA CALIBRATION

77	17.757	2490	2494	2500	rVB3	33069	44013	0.25%	0.030%
78	17.822	2500	2505	2509	rBV	269544	344726	1.94%	0.238%
79	18.098	2547	2552	2557	rBV3	56224	84946	0.48%	0.059%
80	18.487	2609	2618	2627	rBV	1248150	1699255	9.54%	1.174%
81	18.610	2635	2639	2646	rVB8	58693	88047	0.49%	0.061%
82	18.781	2664	2668	2669	rBV4	25219	33546	0.19%	0.023%
83	18.810	2670	2673	2677	rVB2	66810	77663	0.44%	0.054%
84	18.963	2692	2699	2703	rBV2	174668	235135	1.32%	0.163%
85	19.051	2710	2714	2717	rBV2	66039	85463	0.48%	0.059%
86	19.092	2718	2721	2723	rVV3	42548	57388	0.32%	0.040%
87	19.116	2723	2725	2729	rVB	59198	70451	0.40%	0.049%
88	19.475	2780	2786	2794	rBV	5958939	7862622	44.14%	5.434%
89	20.969	3036	3040	3046	rBV	198270	290578	1.63%	0.201%
90	21.239	3081	3086	3092	rBV	3326202	4231007	23.75%	2.924%
91	23.451	3455	3462	3470	rVB2	4427810	8365193	46.96%	5.782%
92	23.598	3481	3487	3499	rVB2	2436007	4678270	26.27%	3.233%

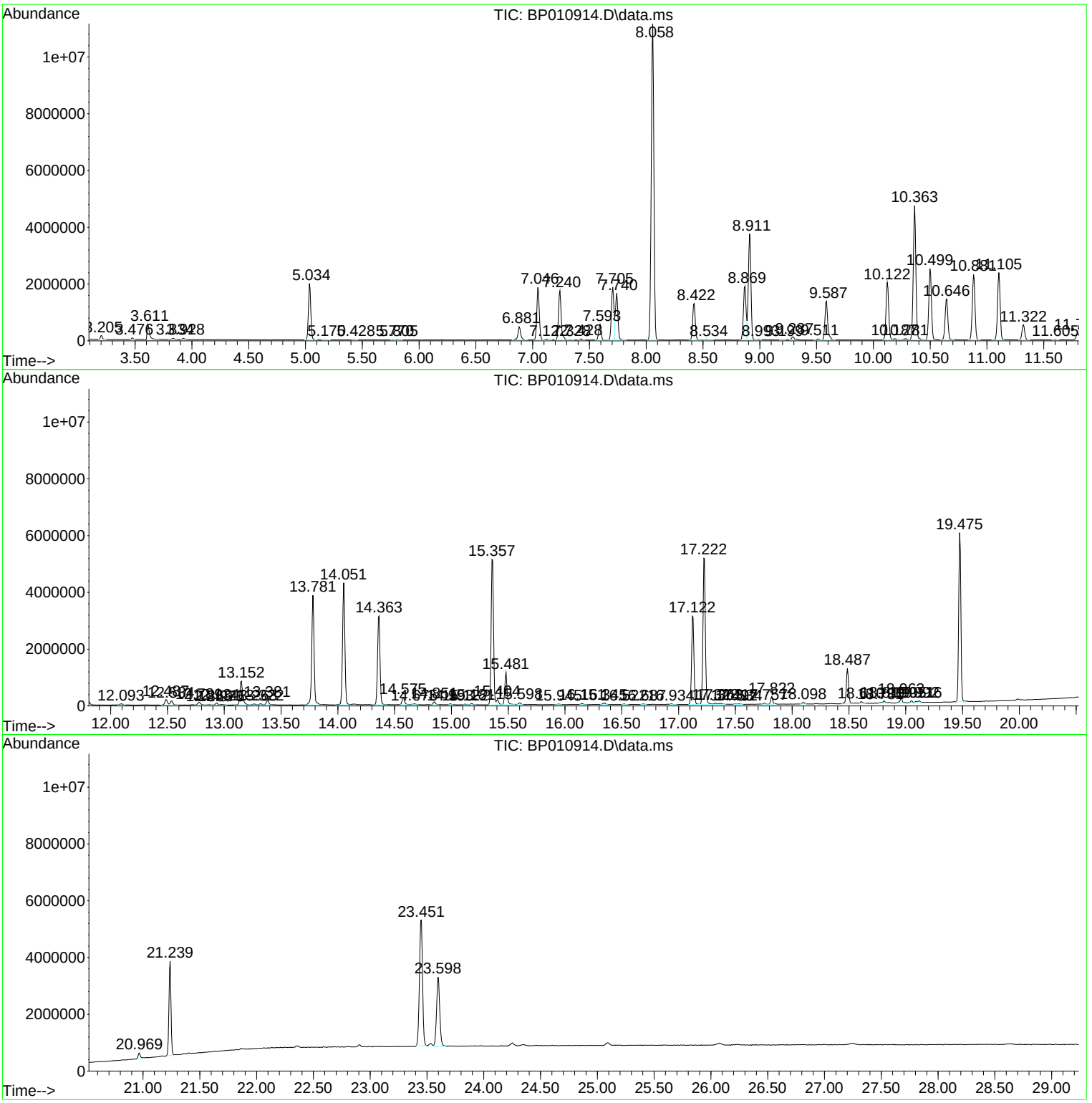
Sum of corrected areas: 144681375

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010914.D
Acq On : 30 Jun 2022 11:30
Operator : CG/JU
Sample : N3523-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JT6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010914.D
Acq On : 30 Jun 2022 11:30
Operator : CG/JU
Sample : N3523-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JT6

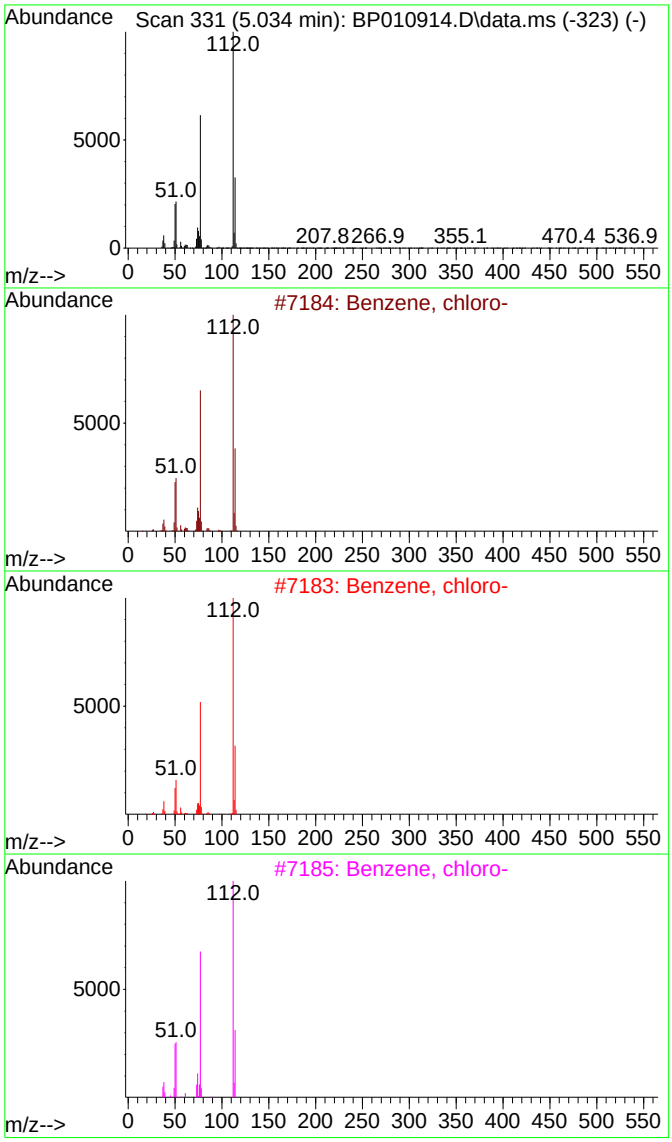
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

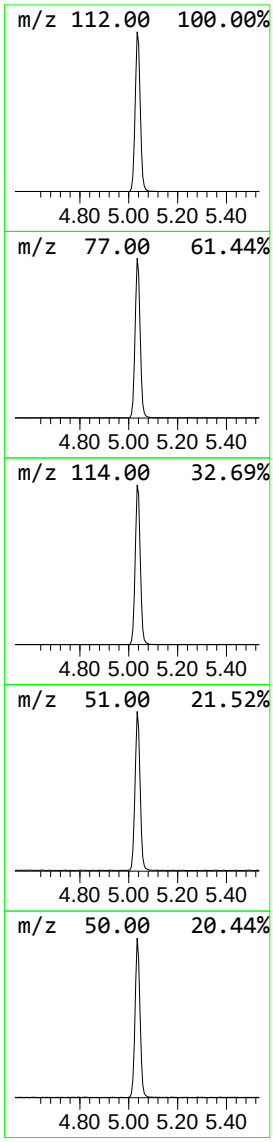
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzene, chloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.034	19.07 ng/ul	2848310	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, chloro-	112	C6H5Cl	000108-90-7	96
2	Benzene, chloro-	112	C6H5Cl	000108-90-7	95
3	Benzene, chloro-	112	C6H5Cl	000108-90-7	95
4	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
5	Benzene, chloro-	112	C6H5Cl	000108-90-7	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010914.D
Acq On : 30 Jun 2022 11:30
Operator : CG/JU
Sample : N3523-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JT6

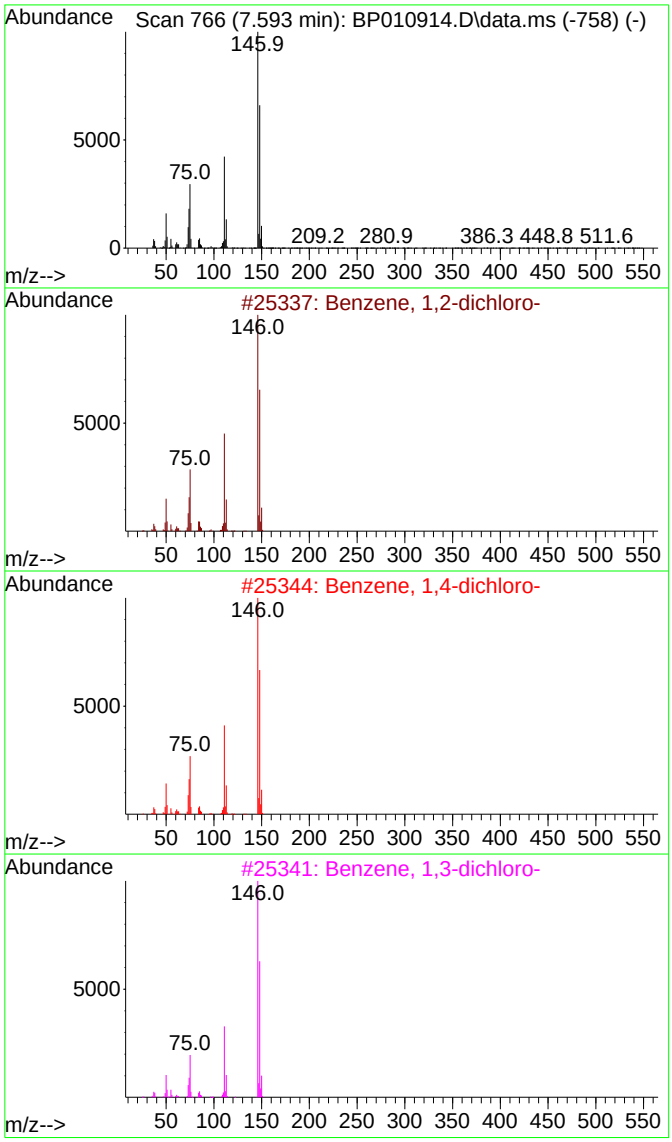
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

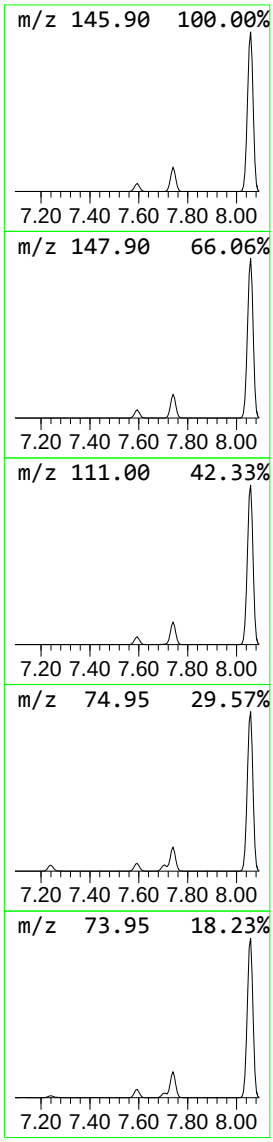
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, 1,2-dichloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.593	5.63 ng/ul	840577	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
2	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
3	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
4	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010914.D
Acq On : 30 Jun 2022 11:30
Operator : CG/JU
Sample : N3523-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JT6

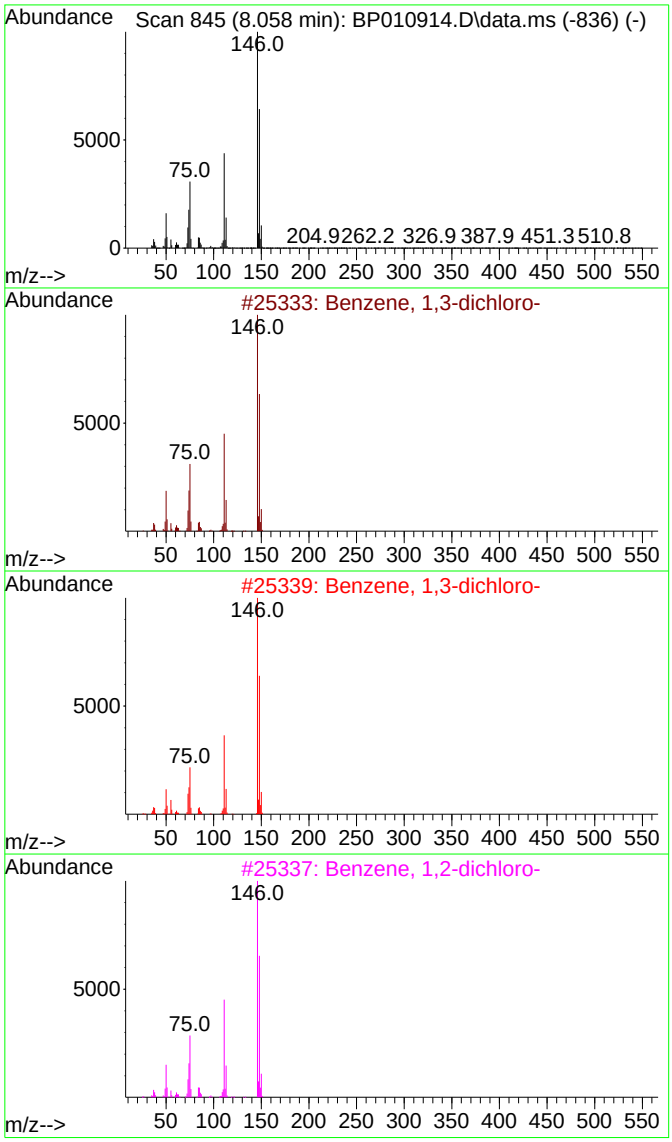
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

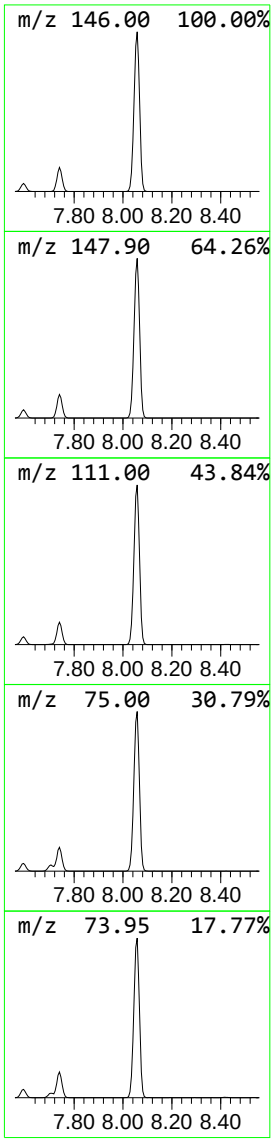
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1,3-dichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.058	119.25 ng/ul	17811600	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
2	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
3	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
4	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
5	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010914.D
Acq On : 30 Jun 2022 11:30
Operator : CG/JU
Sample : N3523-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JT6

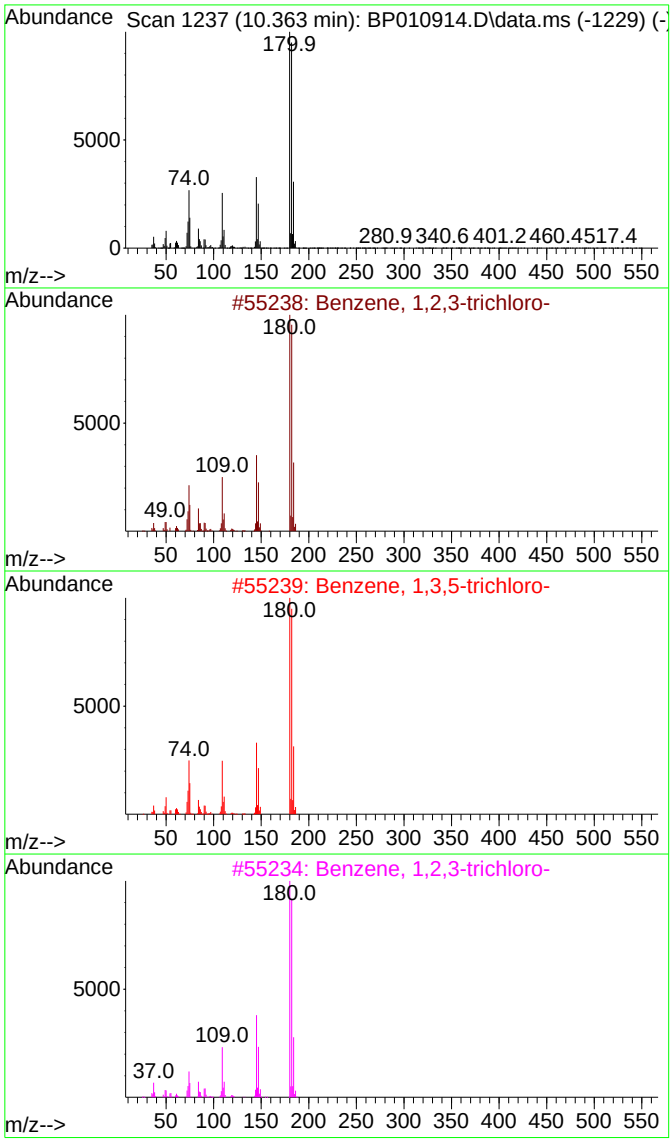
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

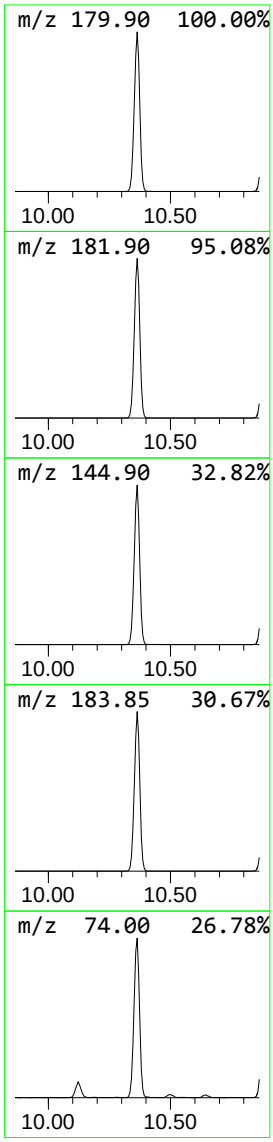
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1,2,3-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.363	37.31 ng/ul	7517940	Naphthalene-d8	10.499

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	99
2	Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
3	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
4	Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	97
5	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010914.D
Acq On : 30 Jun 2022 11:30
Operator : CG/JU
Sample : N3523-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JT6

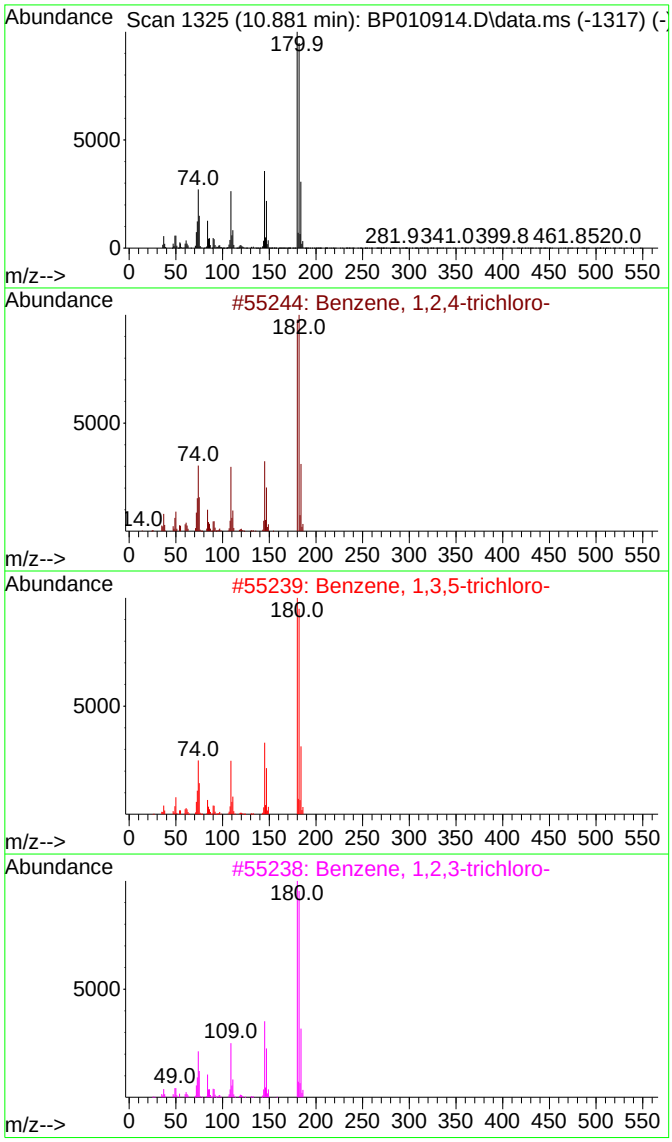
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

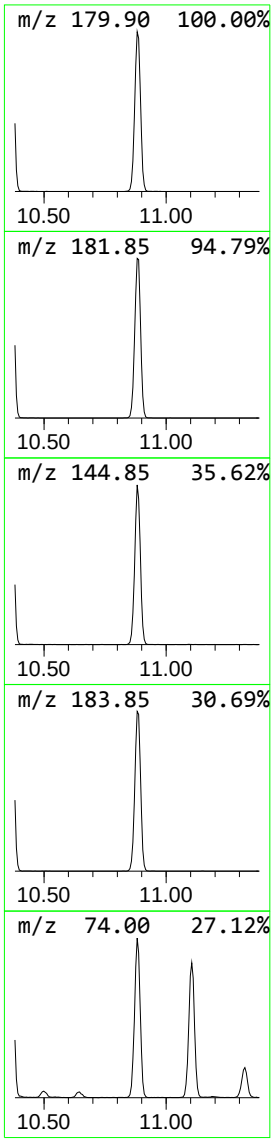
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1,2,4-trichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.881	18.79 ng/ul	3785810	Naphthalene-d8	10.499

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	98
2	Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
3	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
4	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
5	Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	97





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010914.D
Acq On : 30 Jun 2022 11:30
Operator : CG/JU
Sample : N3523-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JT6

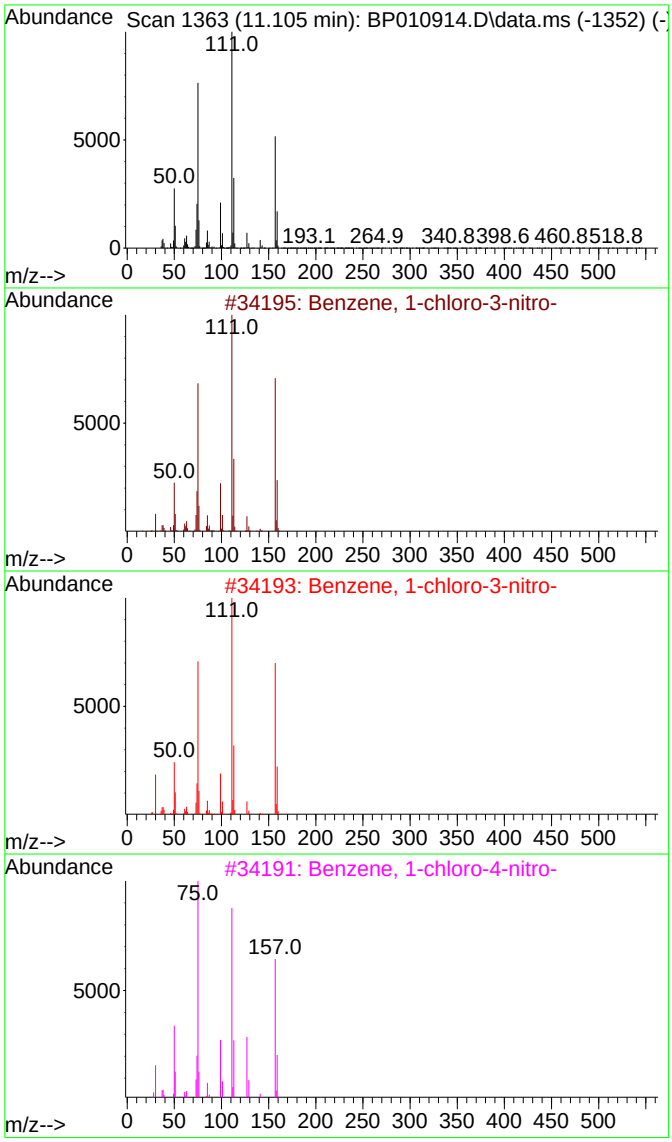
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

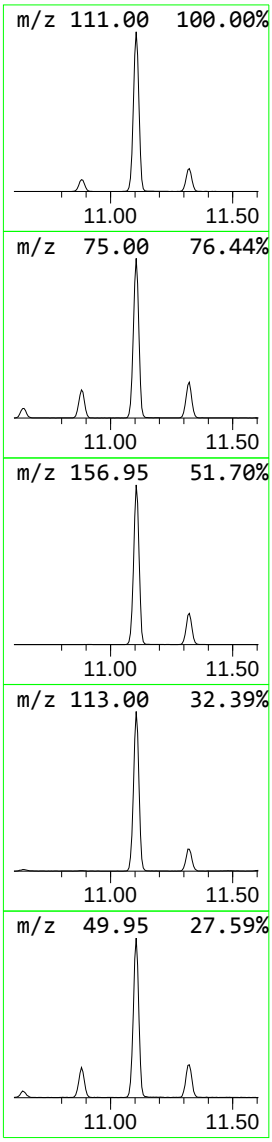
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1-chloro-3-nitro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.105	18.65 ng/ul	3757320	Naphthalene-d8	10.499

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	97
2	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	94
3	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
4	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	92
5	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010914.D
Acq On : 30 Jun 2022 11:30
Operator : CG/JU
Sample : N3523-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JT6

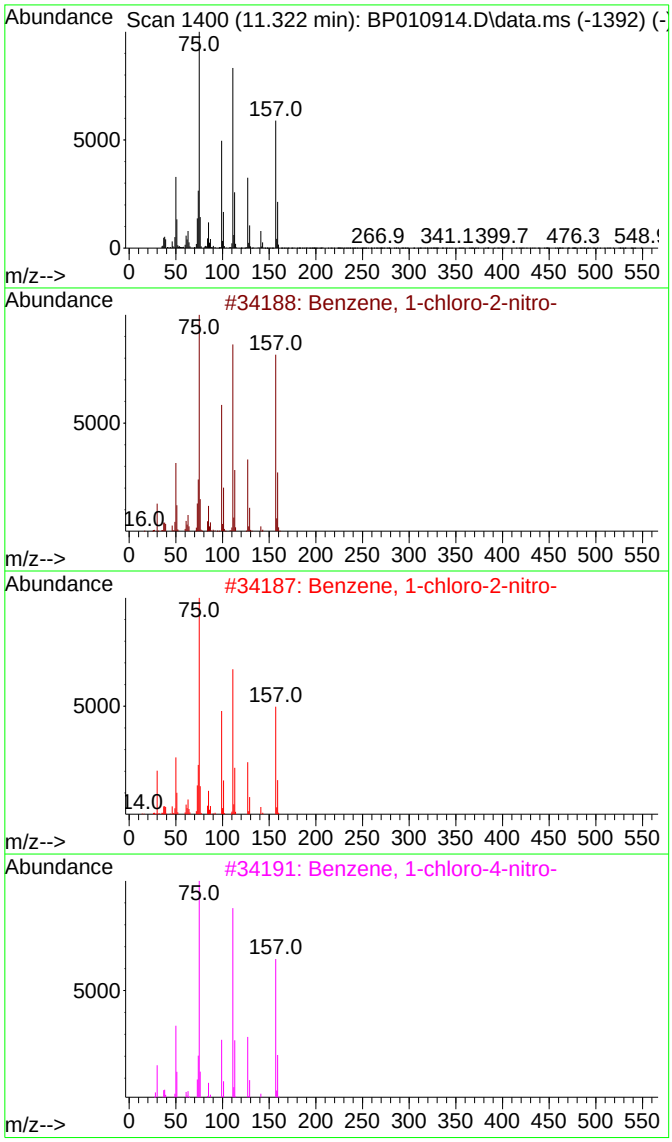
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

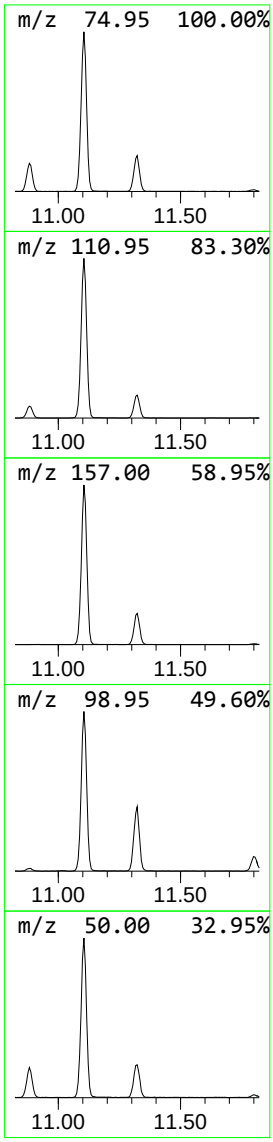
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-chloro-2-nitro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.322	4.45 ng/ul	897265	Naphthalene-d8	10.499

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	98
2	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	97
3	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	97
4	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	96
5	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	96





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010914.D
Acq On : 30 Jun 2022 11:30
Operator : CG/JU
Sample : N3523-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JT6

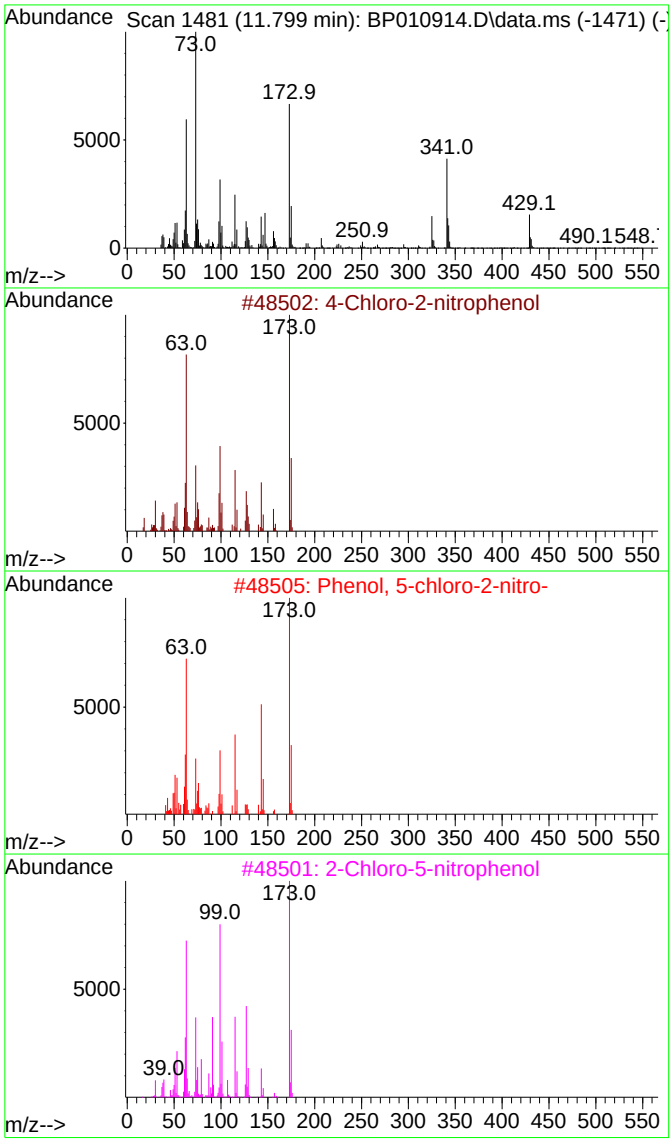
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

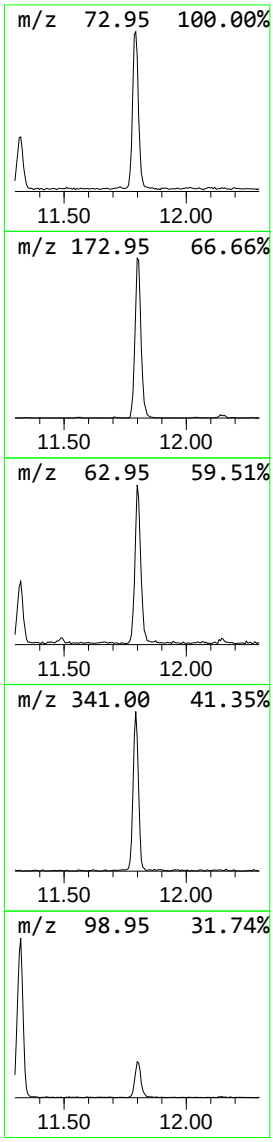
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 4-Chloro-2-nitrophenol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.799	2.20 ng/ul	442597	Naphthalene-d8	10.499

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	98
2	Phenol, 5-chloro-2-nitro-	173	C6H4ClNO3	000611-07-4	62
3	2-Chloro-5-nitrophenol	173	C6H4ClNO3	000619-10-3	62
4	4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	60
5	3-Chloro-6-diazocyclohexa-2,4-di...	154	C6H3ClN2O	080090-95-5	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010914.D
Acq On : 30 Jun 2022 11:30
Operator : CG/JU
Sample : N3523-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JT6

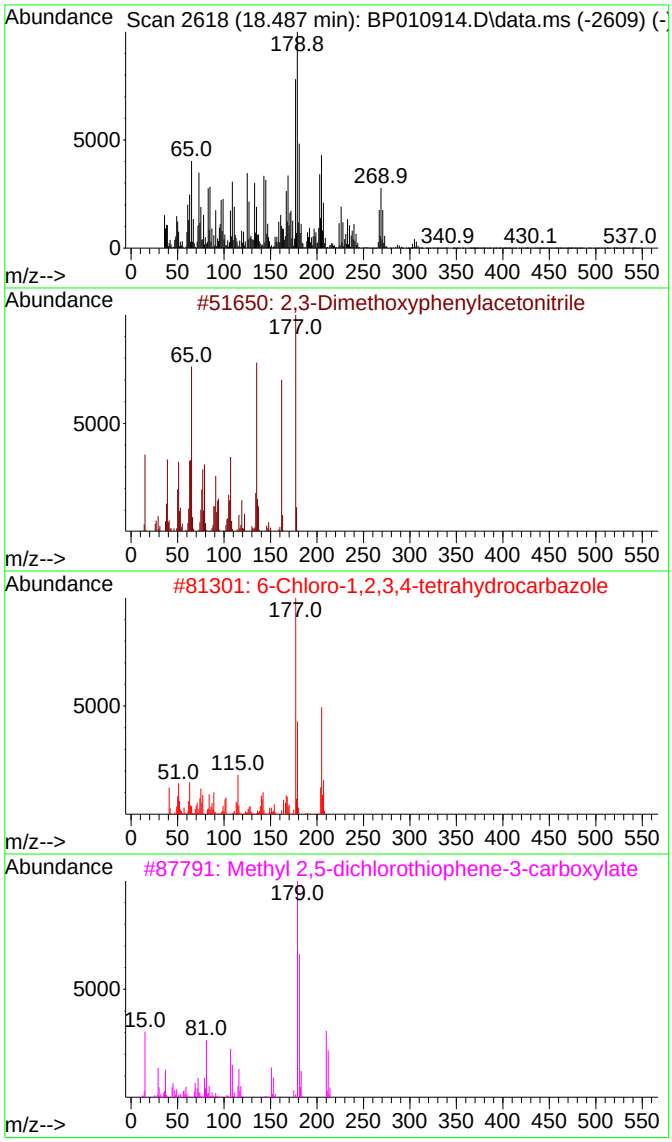
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

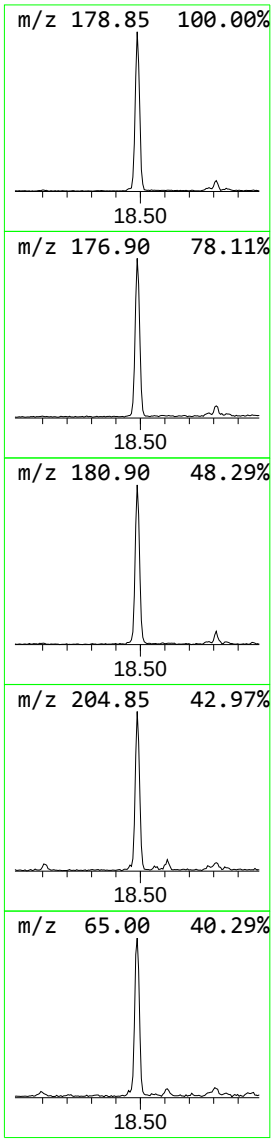
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.486	7.55 ng/ul	1699260	Phenanthrene-d10	17.122

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3-Dimethoxyphenylacetonitrile	177	C10H11NO2	004468-57-9	35
2	6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	35
3	Methyl 2,5-dichlorothiophene-3-c...	210	C6H4Cl2O2S	145129-54-0	14
4	3-(4-Chlorophenyl)-1H-1,2,4-tria...	179	C8H6ClN3	023195-59-7	14
5	Cyclopropane, pentachloro-	212	C3HCl5	006262-51-7	12





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010914.D
Acq On : 30 Jun 2022 11:30
Operator : CG/JU
Sample : N3523-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JT6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA 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
Benzene, chloro-	5.034	19.1	ng/ul	2848310	1	7.705	2987170	20.0
Benzene, 1,2-di...	7.593	5.6	ng/ul	840577	1	7.705	2987170	20.0
Benzene, 1,3-di...	8.058	119.3	ng/ul	17811600	1	7.705	2987170	20.0
Benzene, 1,2,3-...	10.363	37.3	ng/ul	7517940	2	10.499	4030250	20.0
Benzene, 1,2,4-...	10.881	18.8	ng/ul	3785810	2	10.499	4030250	20.0
Benzene, 1-chlo...	11.105	18.6	ng/ul	3757320	2	10.499	4030250	20.0
Benzene, 1-chlo...	11.322	4.5	ng/ul	897265	2	10.499	4030250	20.0
4-Chloro-2-nitr...	11.799	2.2	ng/ul	442597	2	10.499	4030250	20.0
unknown-01	18.486	7.5	ng/ul	1699260	4	17.122	4502000	20.0