LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
Data File : BG053775.D
Acq On : 1 Jun 2022 1:15
Operator : CG/JU
Sample : N3039-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0JP0

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_G\Methods\SFAM-EPA-BG053122.M
Title : SVOA CALIBRATION

Signal : TIC: BG053775.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.187	29	34	42	rBV	47335	100639	11.46%	1.234%
2	3.263	42	47	65	rVB	494392	857203	97.62%	10.515%
3	3.528	88	92	101	rVB4	5487	10276	1.17%	0.126%
4	3.610	104	106	109	rVV3	1098	1183	0.13%	0.015%
5	3.651	111	113	116	rBV4	990	1205	0.14%	0.015%
6	3.951	162	164	174	rVV2	7671	14821	1.69%	0.182%
7	4.021	174	176	179	rVB4	1218	1215	0.14%	0.015%
8	4.186	198	204	210	rVB5	3854	6939	0.79%	0.085%
9	4.662	283	285	288	rVB4	1030	1213	0.14%	0.015%
10	4.832	311	314	315	rBV3	1460	1434	0.16%	0.018%
11	5.408	406	412	422	rVB2	11434	21694	2.47%	0.266%
12	5.555	435	437	441	rVB4	915	1201	0.14%	0.015%
13	5.802	475	479	482	rVV4	1114	1609	0.18%	0.020%
14	6.137	533	536	541	rVB5	971	1831	0.21%	0.022%
15	6.184	541	544	548	rBV5	1363	2049	0.23%	0.025%
16	6.789	646	647	650	rBV3	970	1060	0.12%	0.013%
17	6.853	654	658	659	rVB3	998	1148	0.13%	0.014%
18	7.247	718	725	734	rBV2	19247	37587	4.28%	0.461%
19	7.417	748	754	764	rBV	68373	119019	13.55%	1.460%
20	7.629	783	790	801	rVV2	68255	124450	14.17%	1.527%
21	7.735	805	808	809	rVV2	1224	1351	0.15%	0.017%
22	7.993	845	852	856	rBV4	5154	9745	1.11%	0.120%
23	8.099	864	870	883	rVV2	62957	127503	14.52%	1.564%
24	8.205	885	888	890	rBV3	966	1336	0.15%	0.016%
25	8.375	913	917	921	rBV4	788	1619	0.18%	0.020%
26	8.457	924	931	941	rVV	77955	139591	15.90%	1.712%
27	8.792	981	988	994	rVB	58205	102325	11.65%	1.255%
28	9.256	1058	1067	1071	rBV	79904	160896	18.32%	1.974%
29	9.298	1071	1074	1083	rVV	68864	117707	13.40%	1.444%
30	9.356	1083	1084	1088	rVB2	1267	1078	0.12%	0.013%
31	9.562	1115	1119	1122	rBV2	620	1172	0.13%	0.014%
32	9.709	1138	1144	1152	rBV3	8971	15843	1.80%	0.194%
33	9.979	1184	1190	1200	rVV2	45264	89664	10.21%	1.100%
34	10.467	1270	1273	1275	rVB2	1199	1236	0.14%	0.015%
35	10.520	1275	1282	1291	rBV	92517	173090	19.71%	2.123%
36	10.678	1306	1309	1314	rBV3	889	1596	0.18%	0.020%

LSC Area Percent Report

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Sample : N3039-09
Misc :
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Instrument :
BNA_G
ClientSampleId :
C0JP0

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_G\Methods\SFAM-EPA-BG053122.M
Title : SVOA CALIBRATION

37	10.772	1318	1325	1331	rVB2	58192	107766	12.27%	1.322%
38	10.907	1340	1348	1356	rBV	86032	161907	18.44%	1.986%
39	11.031	1362	1369	1378	rVB	45681	83876	9.55%	1.029%
40	11.295	1408	1414	1421	rBV	37024	64957	7.40%	0.797%
41	11.501	1439	1449	1455	rVV	51036	93676	10.67%	1.149%
42	11.712	1478	1485	1490	rVV2	10095	19720	2.25%	0.242%
43	12.206	1561	1569	1576	rVB3	5938	12294	1.40%	0.151%
44	12.488	1610	1617	1623	rVB2	1834	2819	0.32%	0.035%
45	12.829	1669	1675	1677	rBV3	704	1179	0.13%	0.014%
46	12.864	1677	1681	1685	rVB	3482	4462	0.51%	0.055%
47	12.923	1685	1691	1694	rBV3	2507	3508	0.40%	0.043%
48	13.146	1723	1729	1733	rVB5	2301	3891	0.44%	0.048%
49	13.293	1749	1754	1755	rBV2	2027	1644	0.19%	0.020%
50	13.340	1758	1762	1764	rVB3	1203	1384	0.16%	0.017%
51	13.528	1788	1794	1799	rVV	26365	42886	4.88%	0.526%
52	13.722	1822	1827	1833	rBV5	3964	6949	0.79%	0.085%
53	14.121	1888	1895	1902	rVV	233520	356296	40.57%	4.370%
54	14.339	1930	1932	1938	rVB6	1836	2950	0.34%	0.036%
55	14.415	1938	1945	1953	rBV	234536	380151	43.29%	4.663%
56	14.721	1990	1997	2004	rBV	144772	231765	26.39%	2.843%
57	14.879	2018	2024	2030	rBV2	18673	30195	3.44%	0.370%
58	15.196	2072	2078	2081	rBV4	4005	4939	0.56%	0.061%
59	15.449	2119	2121	2124	rBV2	1061	1111	0.13%	0.014%
60	15.520	2129	2133	2138	rVB2	3703	5311	0.60%	0.065%
61	15.714	2159	2166	2174	rVV	328476	548137	62.42%	6.724%
62	15.813	2177	2183	2189	rVV	31195	48857	5.56%	0.599%
63	15.884	2193	2195	2199	rBV3	969	1159	0.13%	0.014%
64	15.954	2202	2207	2214	rBV3	2901	5168	0.59%	0.063%
65	16.048	2221	2223	2227	rVB3	1034	1176	0.13%	0.014%
66	16.489	2295	2298	2301	rBV3	946	1323	0.15%	0.016%
67	16.689	2330	2332	2335	rVB2	963	1128	0.13%	0.014%
68	16.777	2346	2347	2352	rVB2	1057	1228	0.14%	0.015%
69	16.824	2352	2355	2356	rBV2	1062	1101	0.13%	0.014%
70	17.212	2418	2421	2424	rBV4	1071	1570	0.18%	0.019%
71	17.382	2447	2450	2451	rVB2	1455	1244	0.14%	0.015%
72	17.394	2451	2452	2456	rVB3	1182	1417	0.16%	0.017%
73	17.464	2456	2464	2471	rBV	197942	308766	35.16%	3.787%
74	17.564	2473	2481	2489	rBV	418073	661045	75.28%	8.109%
75	17.694	2501	2503	2505	rBV	1347	1115	0.13%	0.014%
76	17.741	2509	2511	2515	rBV5	1277	1164	0.13%	0.014%

LSC Area Percent Report

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Sample : N3039-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0JP0

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

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.M
Title : SVOA CALIBRATION

77	18.017	2557	2558	2563	rBV3	890	1179	0.13%	0.014%
78	18.052	2563	2564	2567	rVV2	1122	1280	0.15%	0.016%
79	18.075	2567	2568	2569	rVV	1712	1129	0.13%	0.014%
80	18.146	2577	2580	2581	rBV3	1117	1074	0.12%	0.013%
81	18.422	2625	2627	2635	rVB4	3449	5578	0.64%	0.068%
82	18.493	2635	2639	2641	rVB4	1396	1597	0.18%	0.020%
83	18.540	2645	2647	2650	rVB4	1486	1143	0.13%	0.014%
84	18.610	2656	2659	2660	rBV2	1036	1157	0.13%	0.014%
85	18.828	2691	2696	2702	rBV2	71931	101149	11.52%	1.241%
86	18.892	2705	2707	2710	rBV4	1267	1403	0.16%	0.017%
87	18.957	2716	2718	2721	rVB3	1774	1496	0.17%	0.018%
88	19.039	2728	2732	2734	rBV4	2137	3029	0.34%	0.037%
89	19.162	2749	2753	2758	rVB6	3449	5265	0.60%	0.065%
90	19.209	2758	2761	2763	rBV3	1683	1752	0.20%	0.021%
91	19.415	2791	2796	2800	rVB5	6295	9560	1.09%	0.117%
92	19.480	2804	2807	2812	rVB2	5243	7257	0.83%	0.089%
93	19.715	2845	2847	2849	rBV3	1572	1279	0.15%	0.016%
94	19.844	2862	2869	2876	rBV	558748	860905	98.04%	10.560%
95	20.067	2905	2907	2912	rBV3	2624	3413	0.39%	0.042%
96	21.401	3130	3134	3139	rBV8	7300	14291	1.63%	0.175%
97	21.471	3142	3146	3150	rVB2	10921	15492	1.76%	0.190%
98	21.742	3186	3192	3200	rVB	234796	393012	44.75%	4.821%
99	24.791	3702	3711	3723	rVB2	298948	878141	100.00%	10.772%
100	25.014	3741	3749	3760	rVB2	132615	376084	42.83%	4.613%

Sum of corrected areas: 8152427

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
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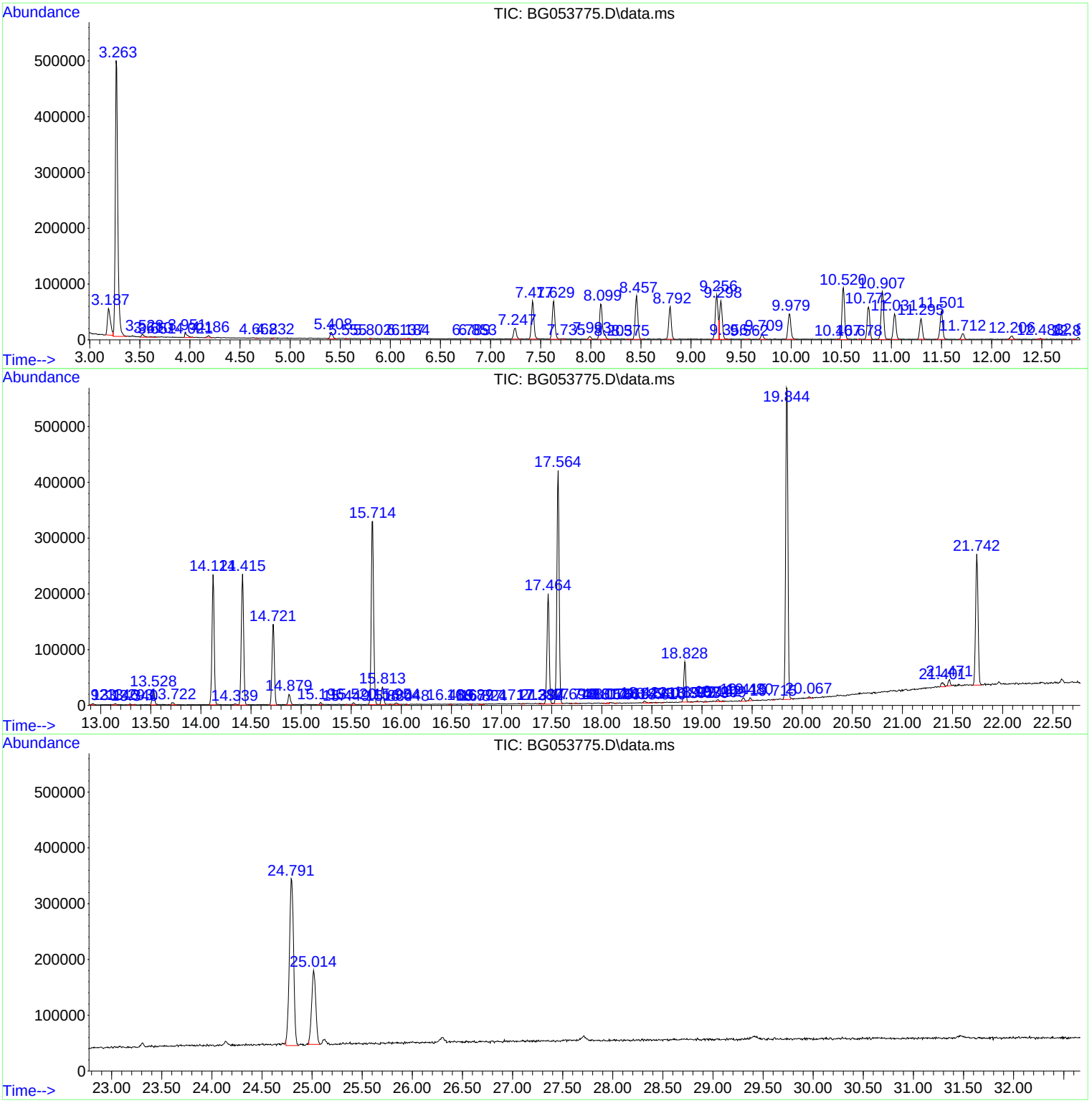
Sample : N3039-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0JP0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
Data File : BG053775.D
Acq On : 1 Jun 2022 1:15
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Sample : N3039-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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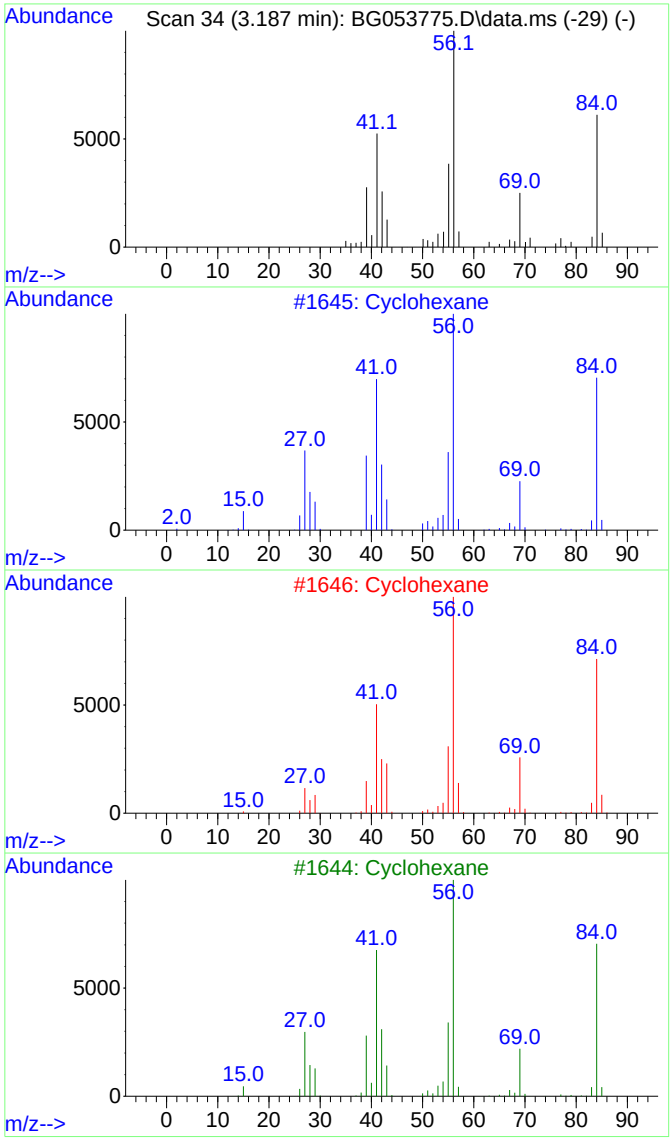
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.M
Quant Title : SVOA CALIBRATION

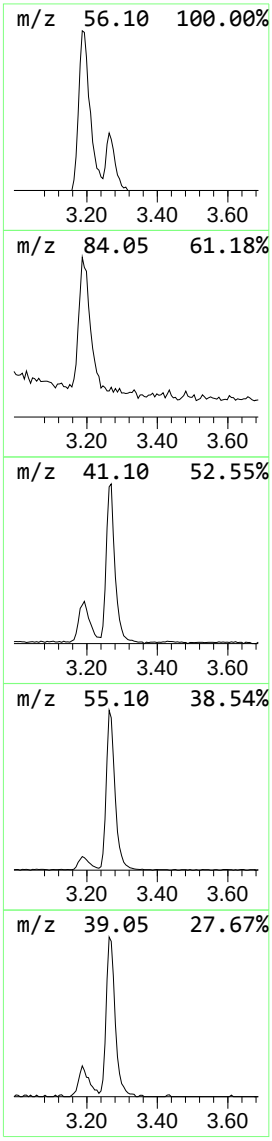
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Cyclic3.187 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.187	15.79 ng/ul	100639	1,4-Dichlorobenzene-d4	8.105

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane	84	C6H12	000110-82-7	90
2	Cyclohexane	84	C6H12	000110-82-7	90
3	Cyclohexane	84	C6H12	000110-82-7	90
4	Cyclohexane	84	C6H12	000110-82-7	87
5	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
Data File : BG053775.D
Acq On : 1 Jun 2022 1:15
Operator : CG/JU
Sample : N3039-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0JP0

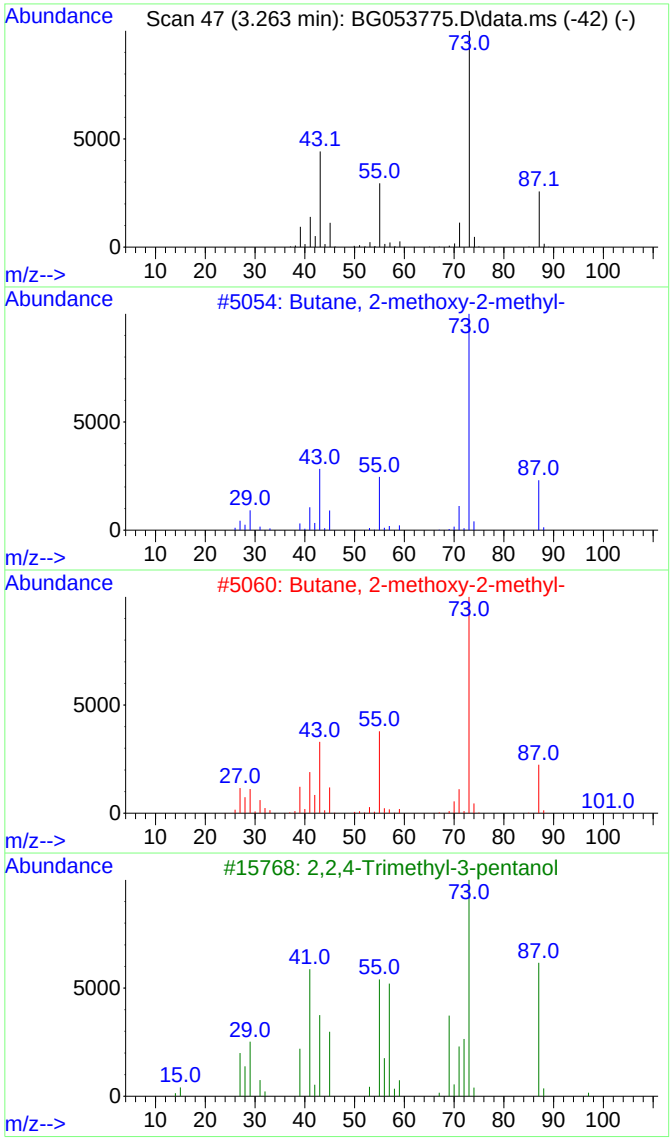
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.M
Quant Title : SVOA CALIBRATION

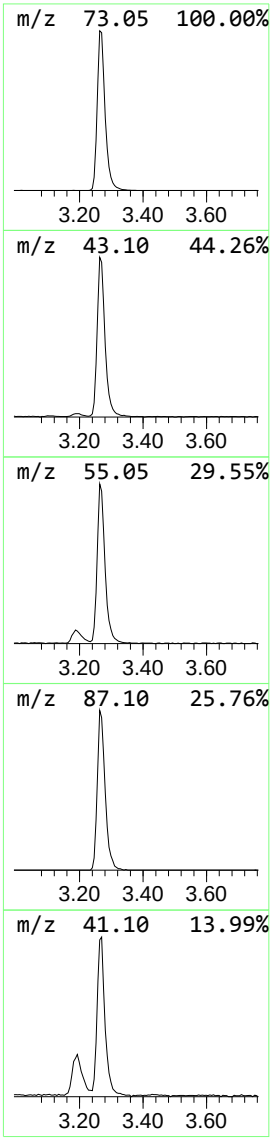
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.263	134.46 ng/ul	857203	1,4-Dichlorobenzene-d4	8.105

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
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Sample : N3039-09
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BNA_G
ClientSampleId :
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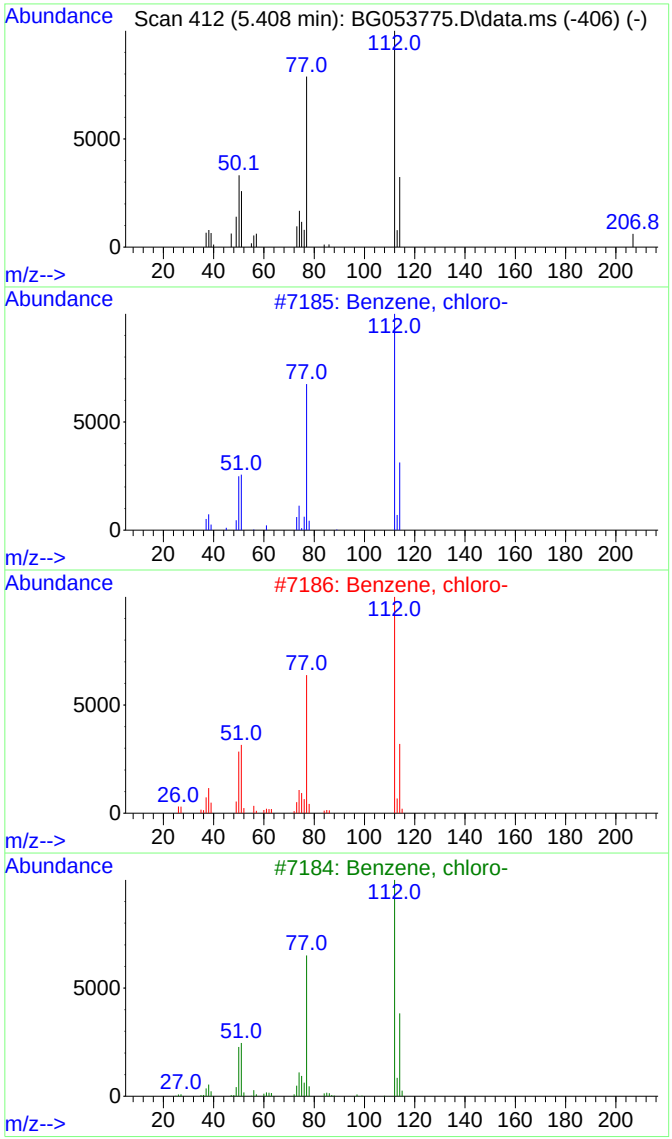
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.M
Quant Title : SVOA CALIBRATION

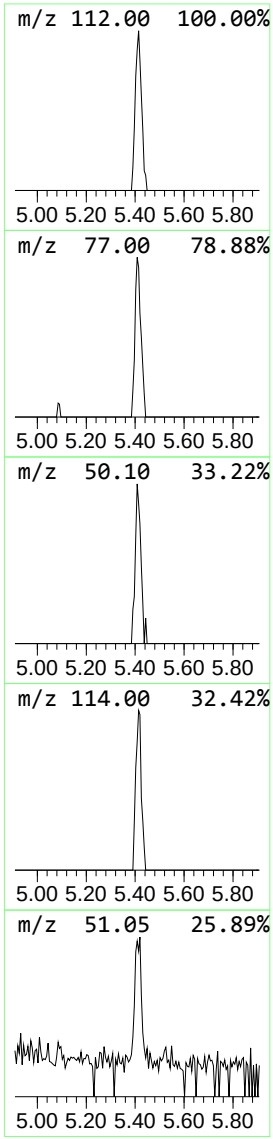
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.408	3.40 ng/ul	21694	1,4-Dichlorobenzene-d4	8.105

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
2	Benzene, chloro-	112	C6H5Cl	000108-90-7	91
3	Benzene, chloro-	112	C6H5Cl	000108-90-7	91
4	Benzene, chloro-	112	C6H5Cl	000108-90-7	80
5	Benzene, chloro-	112	C6H5Cl	000108-90-7	52





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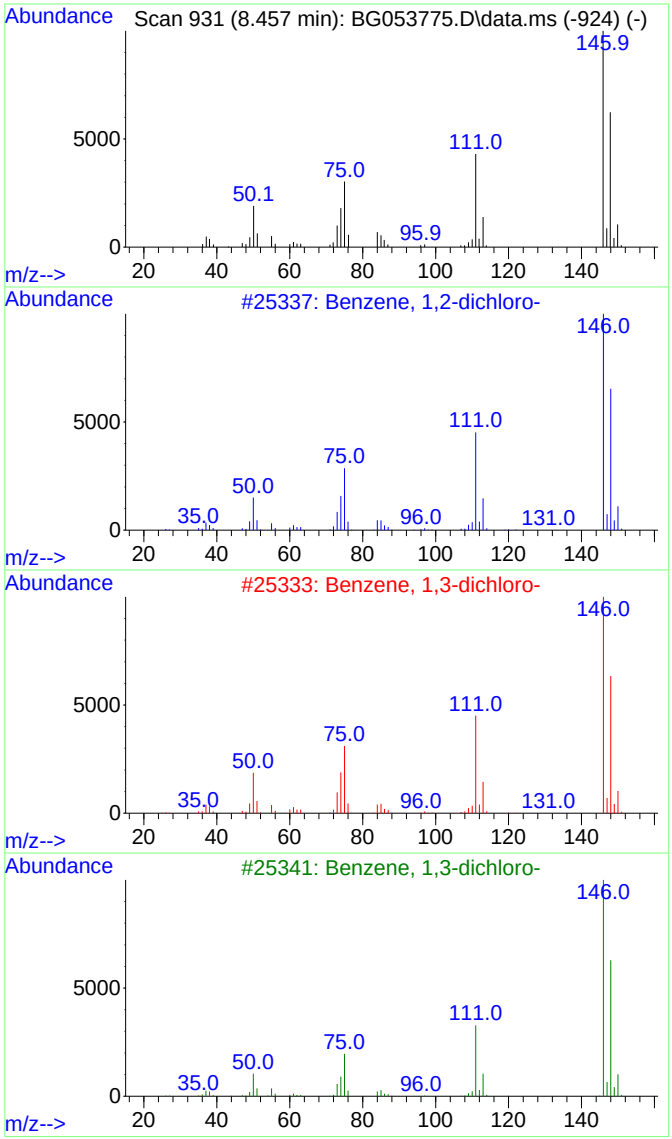
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.M
Quant Title : SVOA CALIBRATION

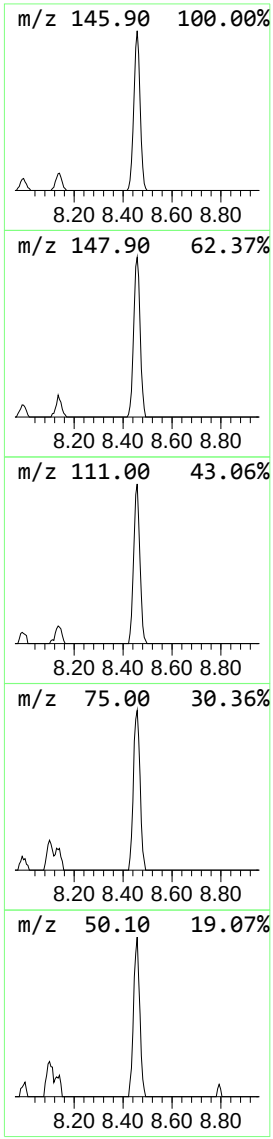
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1,2-dichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.457	21.90 ng/ul	139591	1,4-Dichlorobenzene-d4	8.105

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





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Data File : BG053775.D
Acq On : 1 Jun 2022 1:15
Operator : CG/JU
Sample : N3039-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0JP0

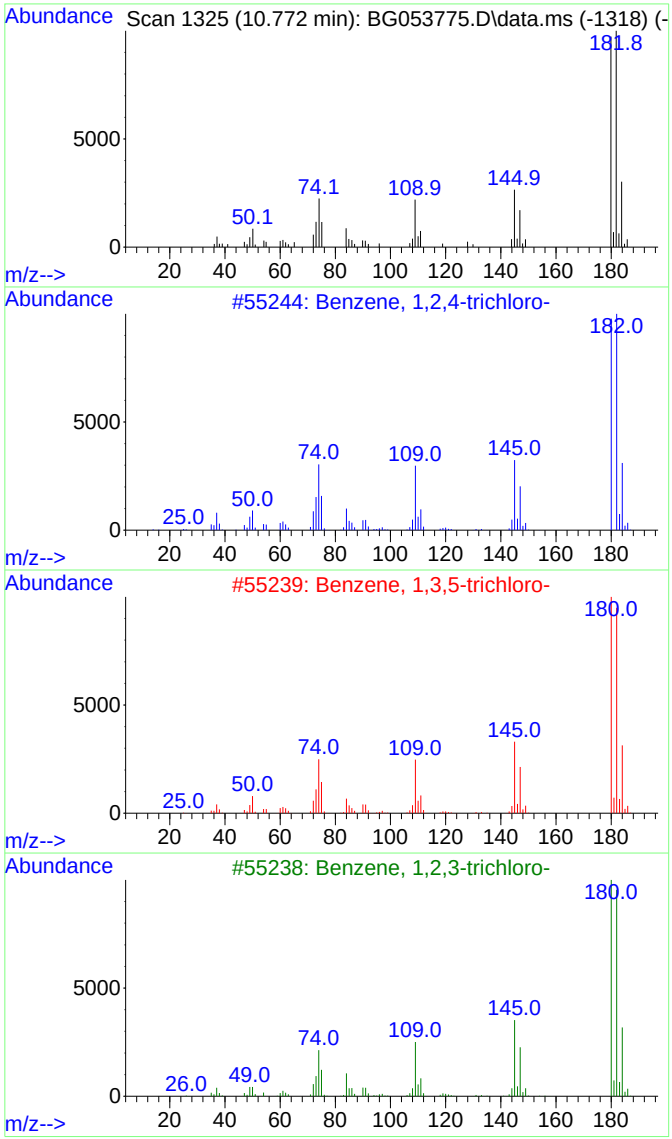
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.M
Quant Title : SVOA CALIBRATION

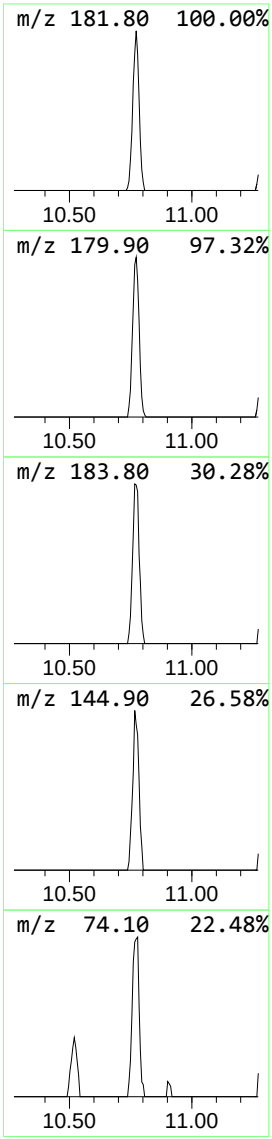
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.772	13.31 ng/ul	107766	Naphthalene-d8	10.907

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, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	96
5	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	96





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Data File : BG053775.D
Acq On : 1 Jun 2022 1:15
Operator : CG/JU
Sample : N3039-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0JP0

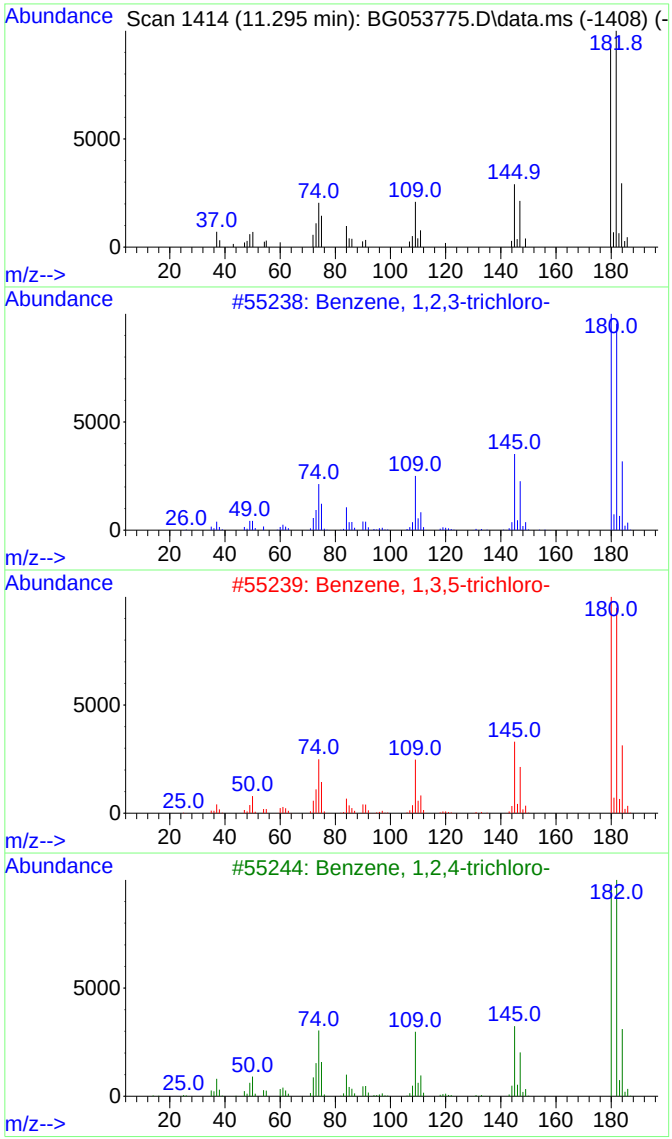
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.M
Quant Title : SVOA CALIBRATION

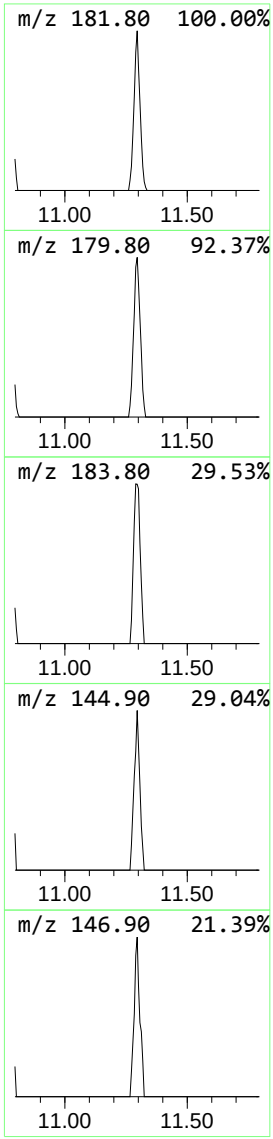
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1,2,3-trichloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.295	8.02 ng/ul	64957	Naphthalene-d8	10.907

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
2	Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
3	Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	96
4	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	96
5	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	96





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
Data File : BG053775.D
Acq On : 1 Jun 2022 1:15
Operator : CG/JU
Sample : N3039-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0JP0

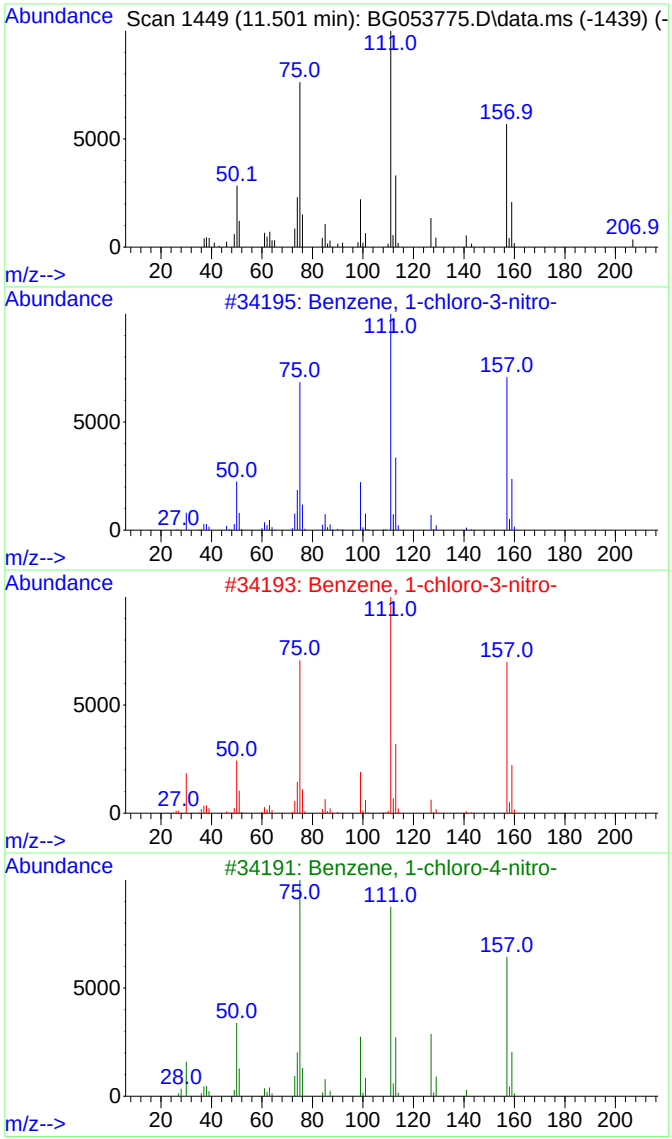
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.M
Quant Title : SVOA CALIBRATION

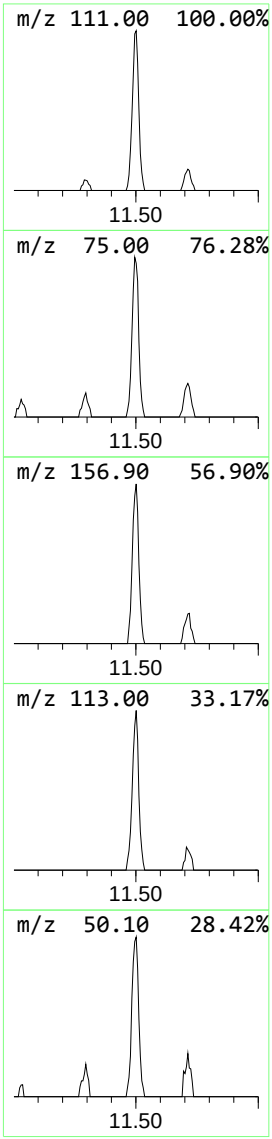
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-chloro-3-nitro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.501	11.57 ng/ul	93676	Naphthalene-d8	10.907

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	96
3	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	93
4	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	91
5	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
Data File : BG053775.D
Acq On : 1 Jun 2022 1:15
Operator : CG/JU
Sample : N3039-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0JP0

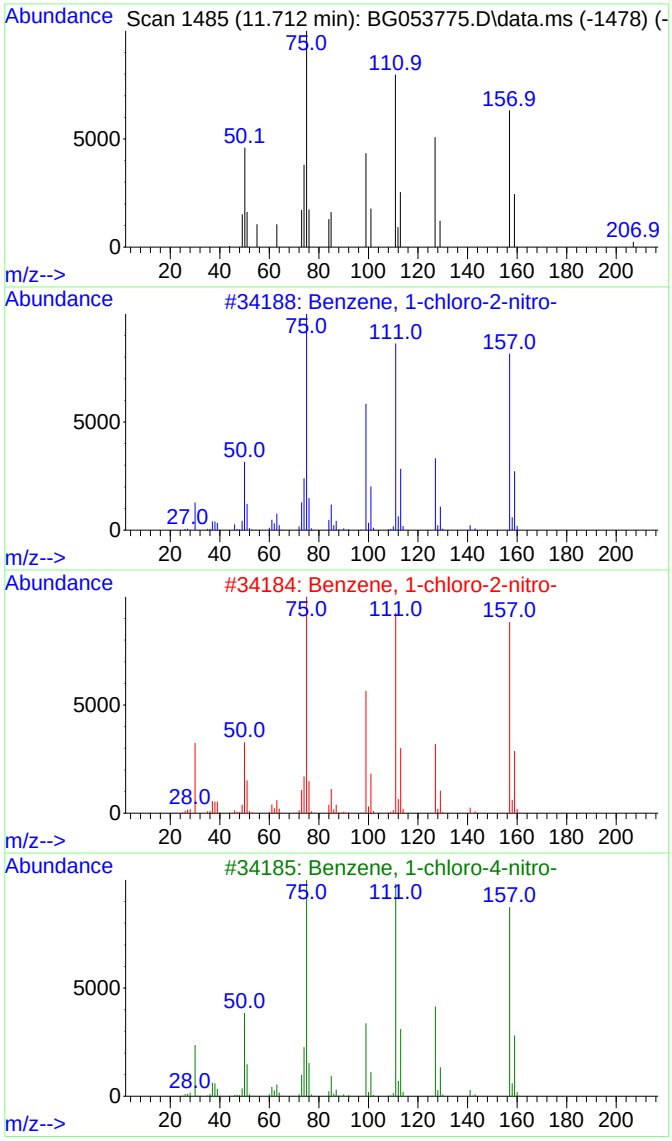
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.M
Quant Title : SVOA CALIBRATION

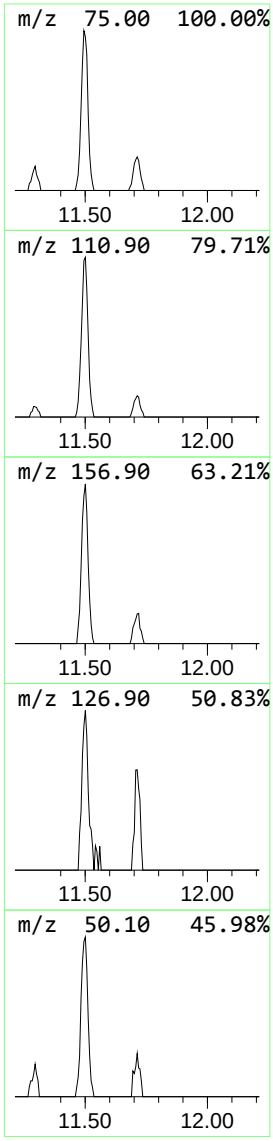
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-chloro-2-nitro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.712	2.44 ng/ul	19720	Naphthalene-d8	10.907

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	91
2	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	91
3	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	90
4	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	90
5	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
Data File : BG053775.D
Acq On : 1 Jun 2022 1:15
Operator : CG/JU
Sample : N3039-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0JP0

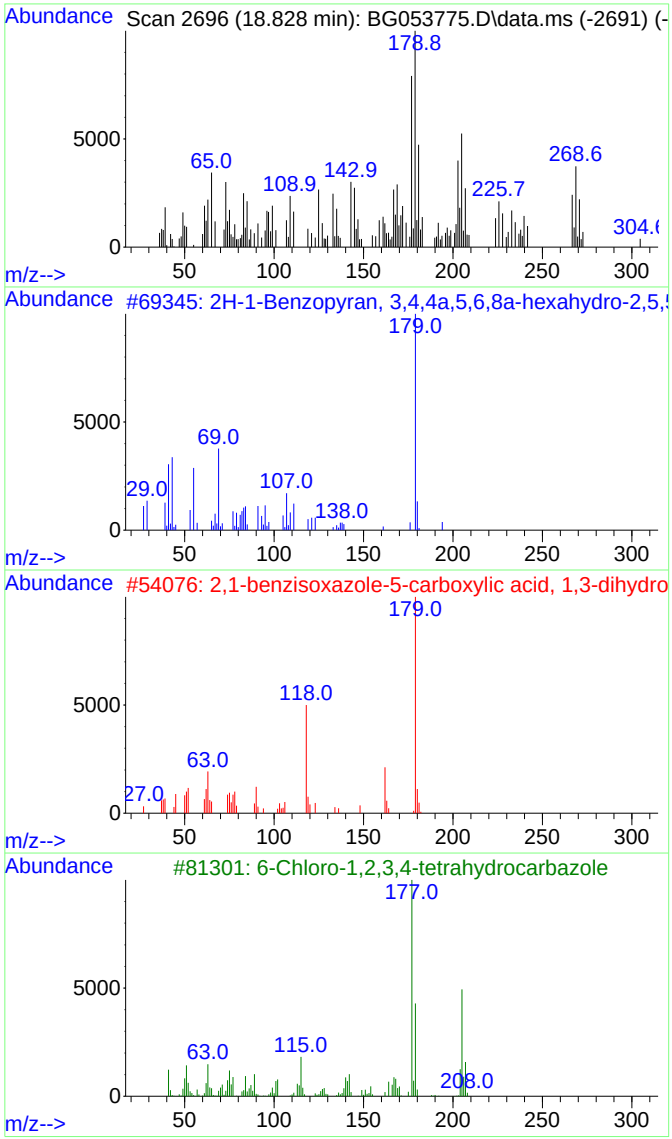
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.M
Quant Title : SVOA CALIBRATION

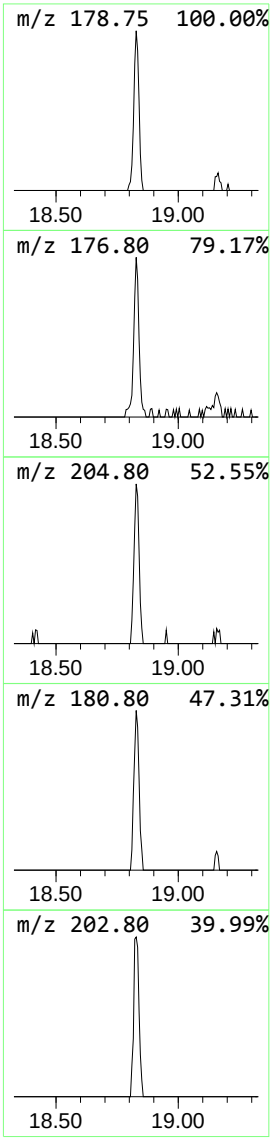
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 2H-1-Benzopyran, 3,4,4a,5,6... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.828	6.55 ng/ul	101149	Phenanthrene-d10	17.464

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-1-Benzopyran, 3,4,4a,5,6,8a-h...	194	C13H22O	041678-32-4	53
2	2,1-benzisoxazole-5-carboxylic a...	179	C8H5NO4	1000404-80-3	35
3	6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	22
4	1-(4-Chlorophenyl)cyclopentaneca...	224	C12H13ClO2	1000484-45-9	20
5	2,4-Dichloro-3-fluoroaniline	179	C6H4Cl2FN	000443-93-6	18





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053122\
Data File : BG053775.D
Acq On : 1 Jun 2022 1:15
Operator : CG/JU
Sample : N3039-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0JP0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG053122.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
(DEL) Alkane: C...	3.187	15.8	ng/ul	100639	1	8.105	127503	20.0
Butane, 2-metho...	3.263	134.5	ng/ul	857203	1	8.105	127503	20.0
Benzene, chloro-	5.408	3.4	ng/ul	21694	1	8.105	127503	20.0
Benzene, 1,2-di...	8.457	21.9	ng/ul	139591	1	8.105	127503	20.0
Benzene, 1,2,4-...	10.772	13.3	ng/ul	107766	2	10.907	161907	20.0
Benzene, 1,2,3-...	11.295	8.0	ng/ul	64957	2	10.907	161907	20.0
Benzene, 1-chlo...	11.501	11.6	ng/ul	93676	2	10.907	161907	20.0
Benzene, 1-chlo...	11.712	2.4	ng/ul	19720	2	10.907	161907	20.0
2H-1-Benzopyran...	18.828	6.5	ng/ul	101149	4	17.464	308766	20.0