LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120222\
Data File : BP012942.D
Acq On : 02 Dec 2022 16:35
Operator : CG/JU
Sample : N5785-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0L75

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-BP120122.M
Title : SVOA CALIBRATION

Signal : TIC: BP012942.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.828	107	109	122	rBV	265884	447235	0.35%	0.062%
2	5.116	322	328	337	rVB	107976	191203	0.15%	0.027%
3	5.310	351	361	367	rBV2	30184492	111290606	86.43%	15.490%
4	6.681	588	594	602	rVB	229552	366252	0.28%	0.051%
5	6.981	640	645	651	rBV	91133	147606	0.11%	0.021%
6	7.163	669	676	678	rBV	360257	613430	0.48%	0.085%
7	7.340	697	706	718	rBV	1349533	2325754	1.81%	0.324%
8	7.540	731	740	752	rBV2	1322018	2706657	2.10%	0.377%
9	7.734	767	773	780	rVB	79688	145107	0.11%	0.020%
10	7.910	793	803	814	rBV	24609628	55496120	43.10%	7.724%
11	8.057	814	828	834	rBV	29087726	104819630	81.40%	14.589%
12	8.387	873	884	892	rBV2	29130623	128770894	100.00%	17.923%
13	8.710	932	939	949	rVB	1117102	1762718	1.37%	0.245%
14	9.175	1010	1018	1021	rBV	1783923	2928555	2.27%	0.408%
15	9.222	1021	1026	1033	rVB	3978644	6634620	5.15%	0.923%
16	9.510	1068	1075	1083	rBV	952671	1559354	1.21%	0.217%
17	9.740	1108	1114	1118	rBV	230137	368633	0.29%	0.051%
18	9.787	1118	1122	1124	rVV	206675	330005	0.26%	0.046%
19	9.828	1124	1129	1135	rVV	512861	954417	0.74%	0.133%
20	9.898	1135	1141	1148	rVV	1478345	2475375	1.92%	0.345%
21	9.957	1148	1151	1160	rVB	132832	216166	0.17%	0.030%
22	10.198	1185	1192	1201	rVB	74705	149350	0.12%	0.021%
23	10.440	1222	1233	1241	rBV	2133323	3710943	2.88%	0.516%
24	10.704	1267	1278	1280	rBV2	30568633	78236648	60.76%	10.889%
25	10.834	1293	1300	1306	rBV	2524273	3995456	3.10%	0.556%
26	10.957	1313	1321	1332	rBV	985247	1668335	1.30%	0.232%
27	11.216	1356	1365	1383	rVB	23315263	46423361	36.05%	6.461%
28	11.422	1392	1400	1407	rBV	2175718	3541571	2.75%	0.493%
29	11.634	1424	1436	1443	rBV	509840	892018	0.69%	0.124%
30	12.104	1502	1516	1528	rBV3	206593	506905	0.39%	0.071%
31	12.416	1552	1569	1578	rBV2	221829	510531	0.40%	0.071%
32	12.840	1621	1641	1649	rBV	3883284	6895650	5.35%	0.960%
33	13.069	1673	1680	1687	rBV2	87391	164252	0.13%	0.023%
34	13.445	1736	1744	1753	rBV	10623360	16419333	12.75%	2.285%
35	13.539	1754	1760	1766	rVV	368852	591286	0.46%	0.082%
36	13.651	1774	1779	1785	rVB2	190883	311606	0.24%	0.043%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120222\
Data File : BP012942.D
Acq On : 02 Dec 2022 16:35
Operator : CG/JU
Sample : N5785-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0L75

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-BP120122.M
Title : SVOA CALIBRATION

37	14.045	1837	1846	1852	rBV	4817571	6853287	5.32%	0.954%
38	14.334	1888	1895	1903	rBV	5608292	8274478	6.43%	1.152%
39	14.639	1940	1947	1957	rBV	4321740	6536385	5.08%	0.910%
40	14.828	1971	1979	1990	rBV	412955	729815	0.57%	0.102%
41	14.957	1996	2001	2012	rVB	452010	659730	0.51%	0.092%
42	15.122	2024	2029	2036	rBV	114073	182782	0.14%	0.025%
43	15.633	2109	2116	2127	rBV	7655817	11150509	8.66%	1.552%
44	15.745	2129	2135	2150	rVV	2429700	3440909	2.67%	0.479%
45	15.998	2167	2178	2191	rVB	560135	851494	0.66%	0.119%
46	16.563	2270	2274	2278	rBV	92395	129403	0.10%	0.018%
47	16.610	2278	2282	2288	rVB5	118699	173889	0.14%	0.024%
48	16.869	2318	2326	2332	rBV	188298	292011	0.23%	0.041%
49	17.398	2409	2416	2426	rBV	6089404	8785266	6.82%	1.223%
50	17.498	2426	2433	2444	rVV	9170253	13076622	10.15%	1.820%
51	17.627	2452	2455	2467	rVB6	54023	134898	0.10%	0.019%
52	18.263	2558	2563	2573	rBV	467165	672012	0.52%	0.094%
53	18.345	2573	2577	2583	rVB2	150634	234250	0.18%	0.033%
54	18.733	2635	2643	2652	rBV	1889476	2572798	2.00%	0.358%
55	18.857	2658	2664	2670	rVB8	106008	204092	0.16%	0.028%
56	19.210	2720	2724	2731	rVB	203911	258799	0.20%	0.036%
57	19.304	2735	2740	2744	rBV	127588	191153	0.15%	0.027%
58	19.369	2747	2751	2755	rVB	127344	192290	0.15%	0.027%
59	19.492	2768	2772	2783	rBV	281914	451658	0.35%	0.063%
60	19.727	2805	2812	2825	rBV	12046197	16749558	13.01%	2.331%
61	21.168	3054	3057	3062	rBV	345900	432058	0.34%	0.060%
62	21.480	3105	3110	3118	rBV	8151184	11369330	8.83%	1.582%
63	21.574	3122	3126	3131	rBV	1028756	1270589	0.99%	0.177%
64	22.621	3301	3304	3309	rVB	297070	393884	0.31%	0.055%
65	23.204	3399	3403	3409	rVB	398216	627500	0.49%	0.087%
66	23.827	3501	3509	3523	rBV2	8457519	18431028	14.31%	2.565%
67	23.992	3530	3537	3549	rVB2	5322887	11254151	8.74%	1.566%
68	24.274	3579	3585	3593	rVB	1045139	2127134	1.65%	0.296%
69	24.633	3641	3646	3657	rVB	587889	1209005	0.94%	0.168%

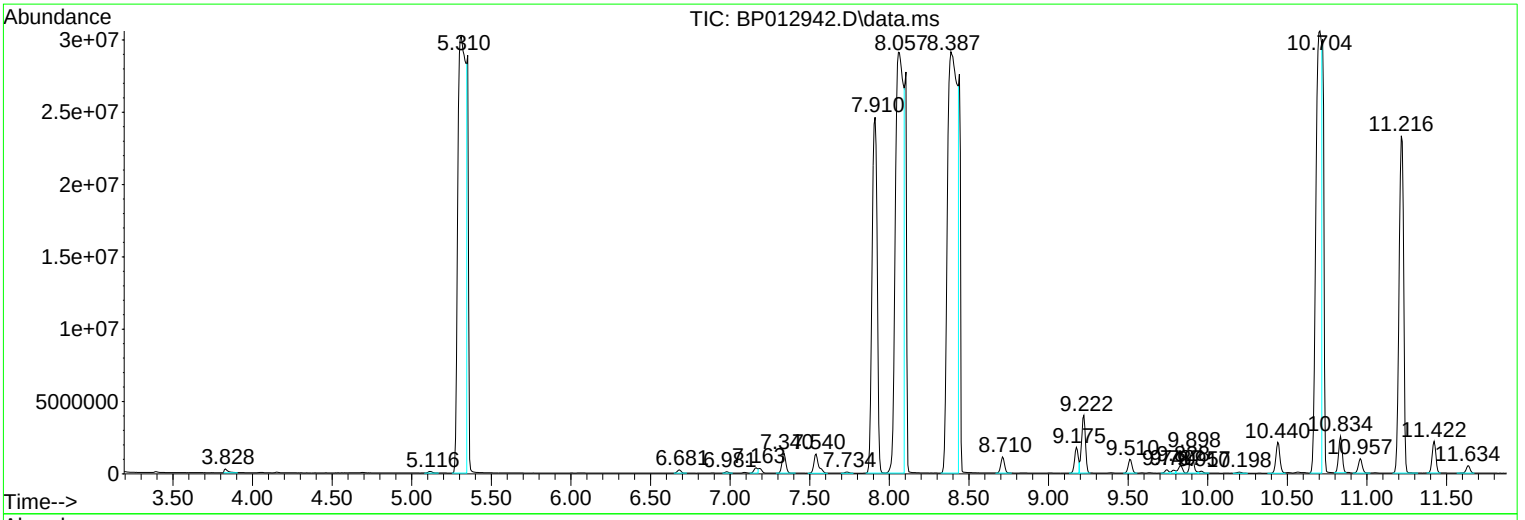
Sum of corrected areas: 718480349

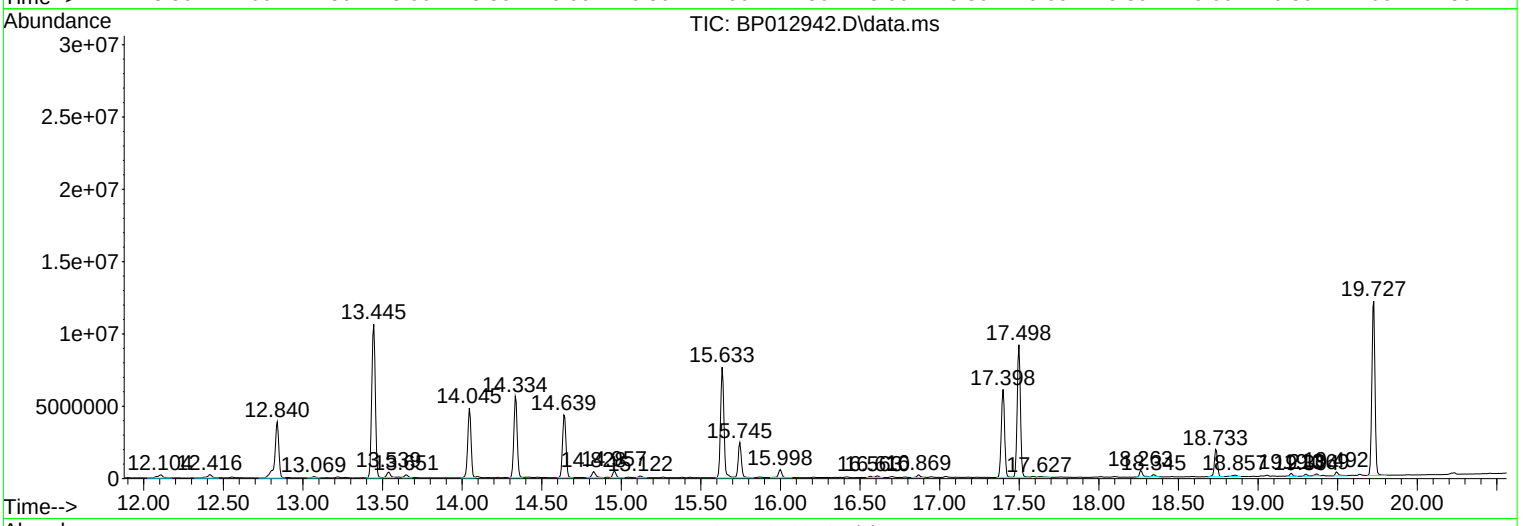
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120222\
Data File : BP012942.D
Acq On : 02 Dec 2022 16:35
Operator : CG/JU
Sample : N5785-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

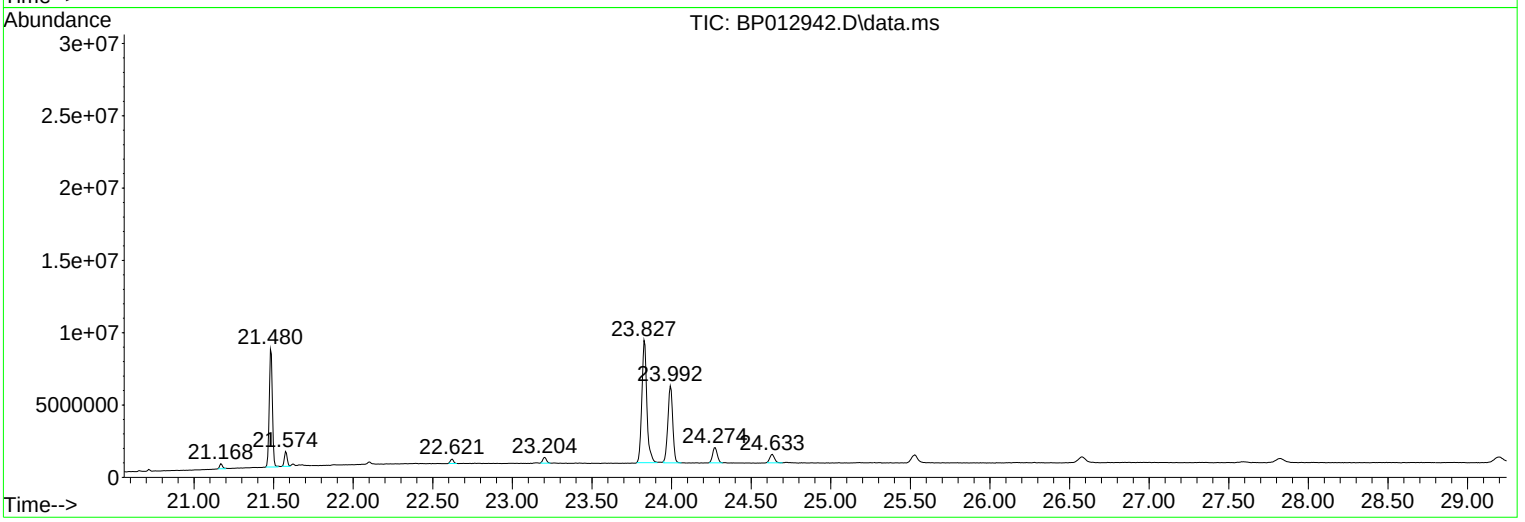
Instrument :
BNA_P
ClientSampleId :
C0L75

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

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







Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120222\
Data File : BP012942.D
Acq On : 02 Dec 2022 16:35
Operator : CG/JU
Sample : N5785-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L75

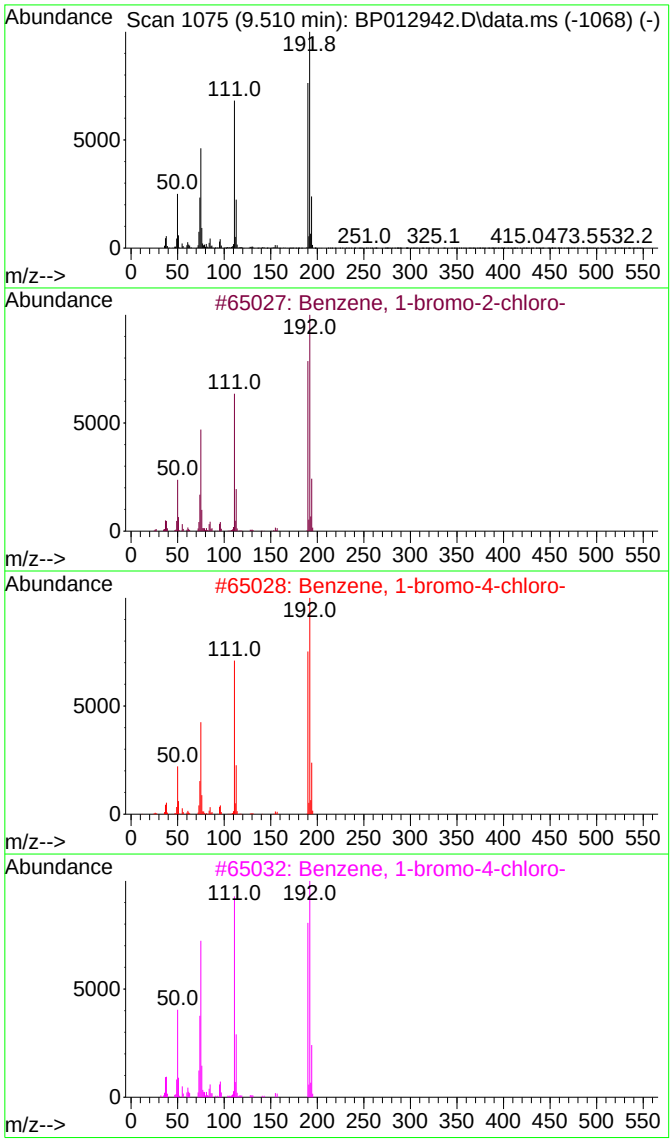
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

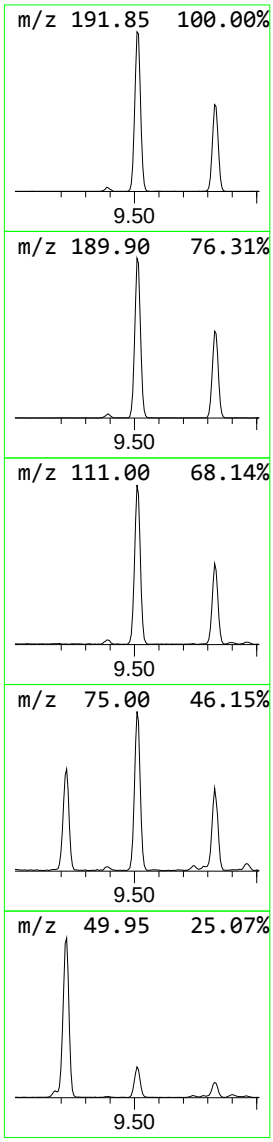
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1-bromo-2-chloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.510	7.81 ng/ul	1559350	Naphthalene-d8	10.834

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	99
2	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
3	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
4	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	97
5	4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	96





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120222\
Data File : BP012942.D
Acq On : 02 Dec 2022 16:35
Operator : CG/JU
Sample : N5785-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L75

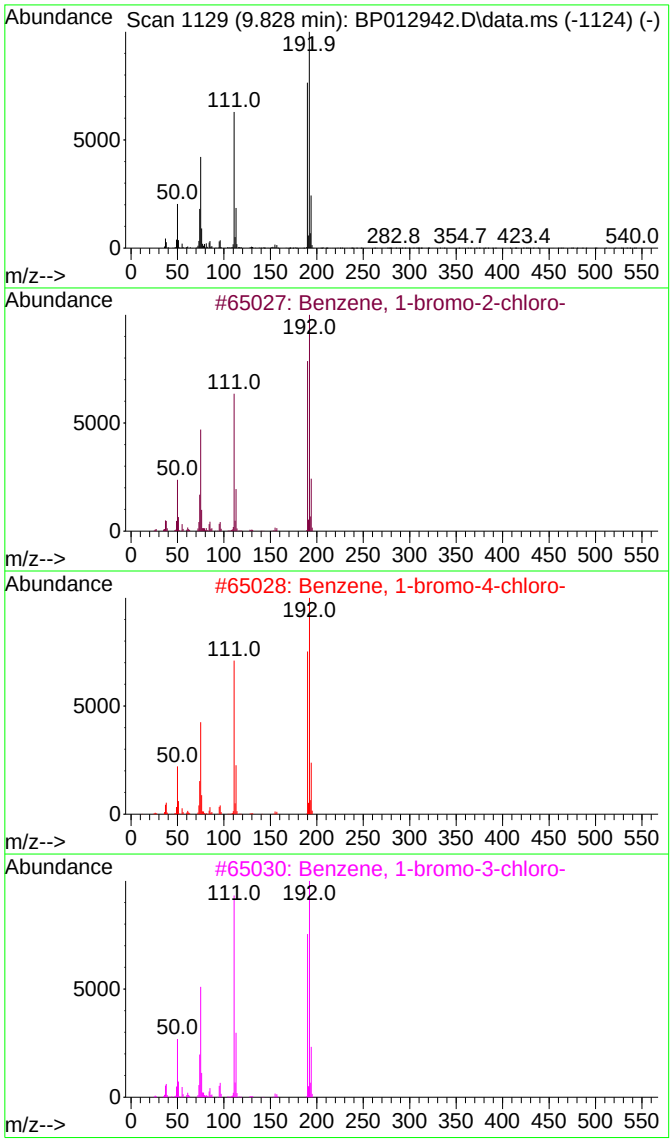
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

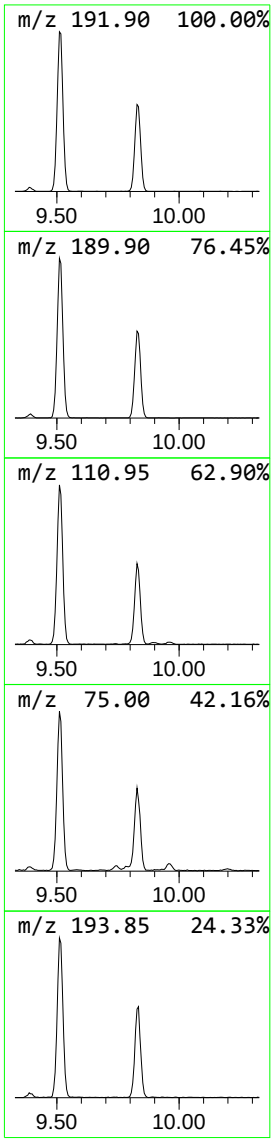
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1-bromo-4-chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.828	4.78 ng/ul	954417	Naphthalene-d8	10.834

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	99
2	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
3	Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97
4	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	95
5	Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120222\
Data File : BP012942.D
Acq On : 02 Dec 2022 16:35
Operator : CG/JU
Sample : N5785-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L75

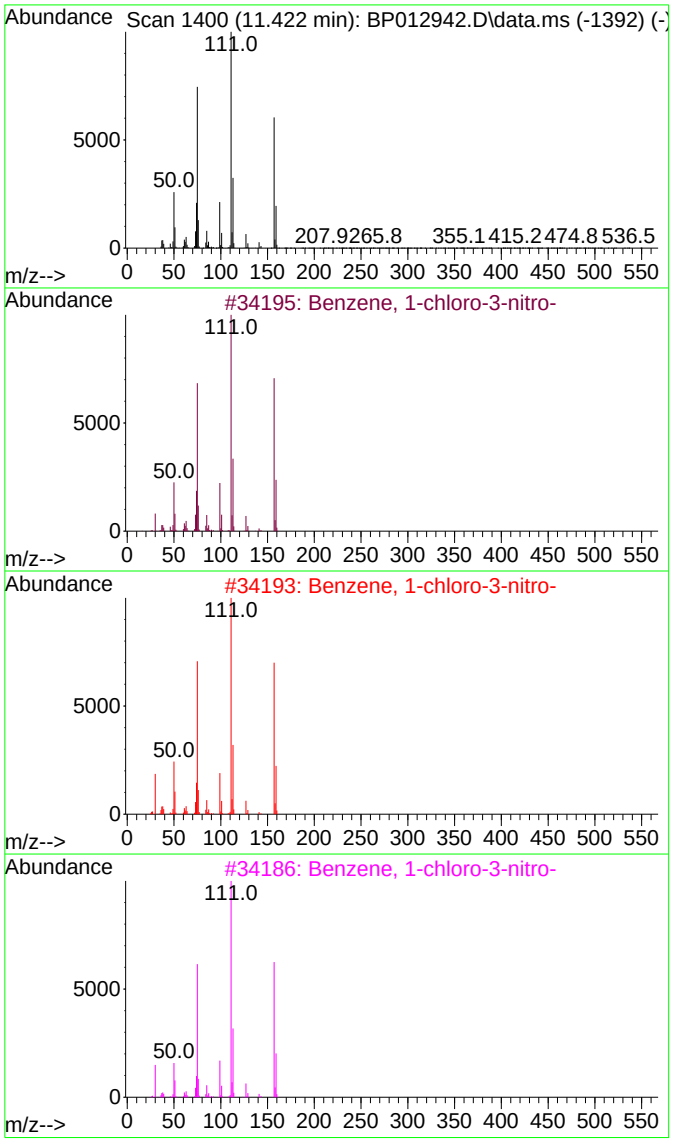
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

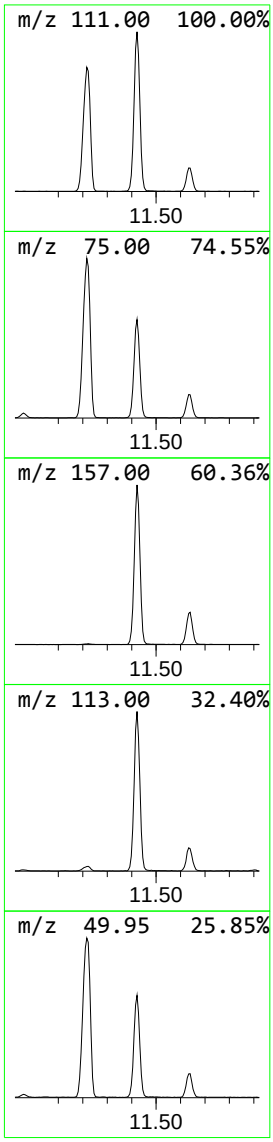
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-chloro-3-nitro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.422	17.73 ng/ul	3541570	Naphthalene-d8	10.834

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	98
2	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	97
3	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	96
4	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
5	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120222\
Data File : BP012942.D
Acq On : 02 Dec 2022 16:35
Operator : CG/JU
Sample : N5785-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L75

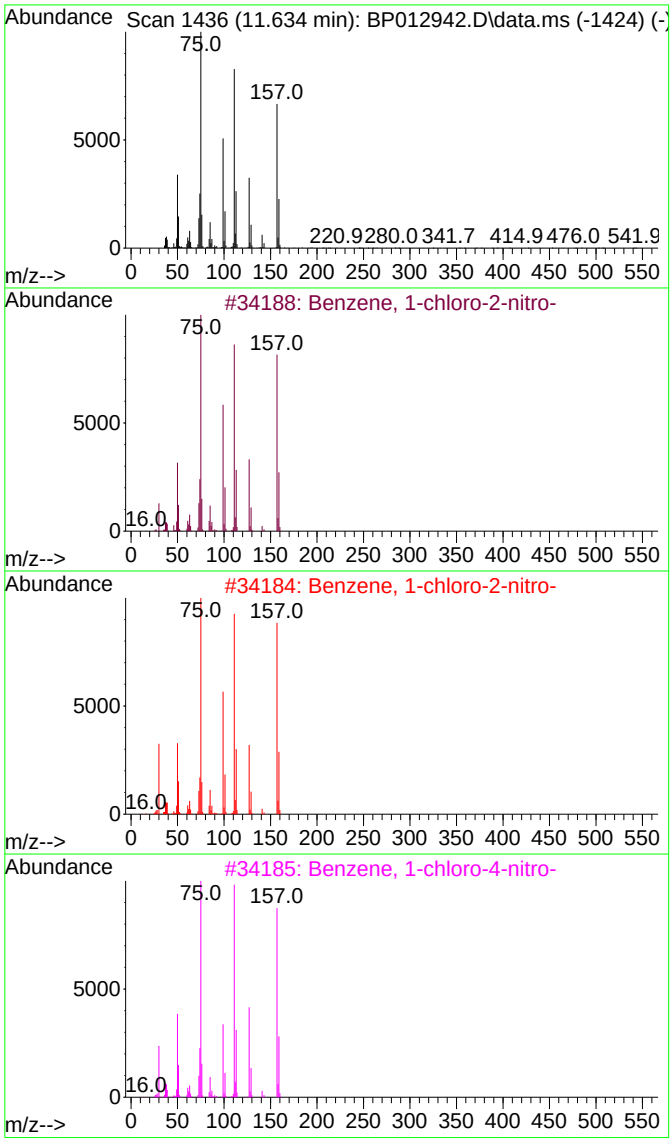
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

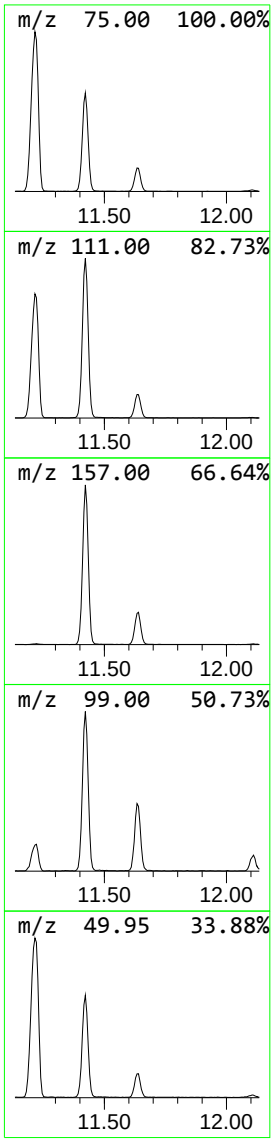
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-chloro-2-nitro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.634	4.47 ng/ul	892018	Naphthalene-d8	10.834

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	99
2	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	98
3	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	98
4	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	97
5	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	97





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120222\
Data File : BP012942.D
Acq On : 02 Dec 2022 16:35
Operator : CG/JU
Sample : N5785-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L75

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

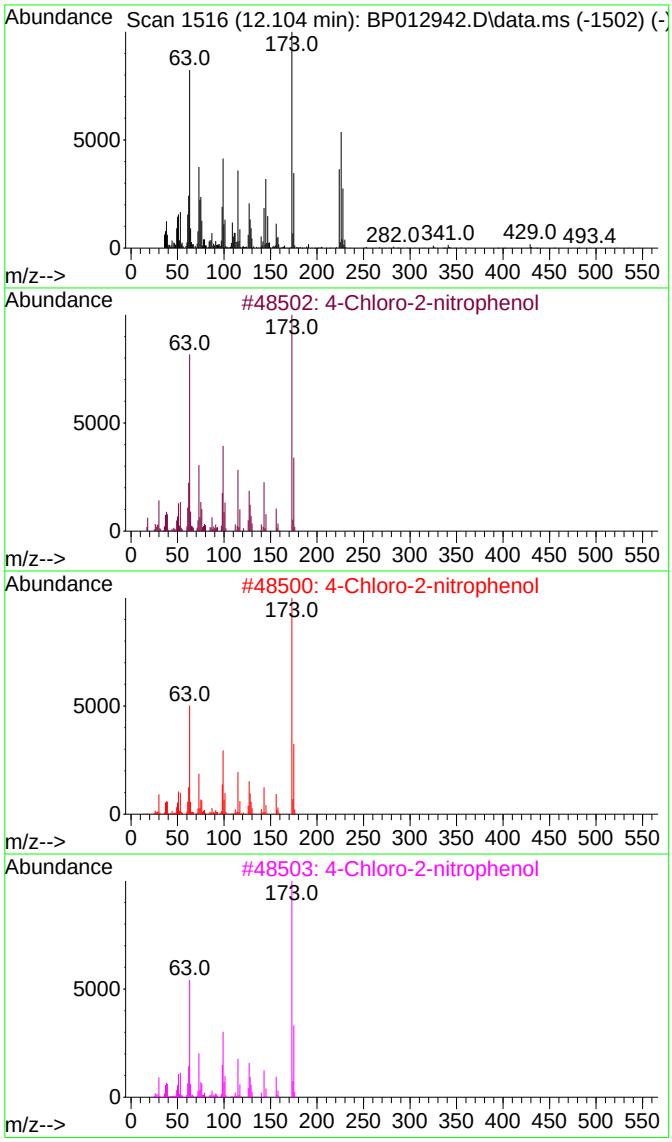
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

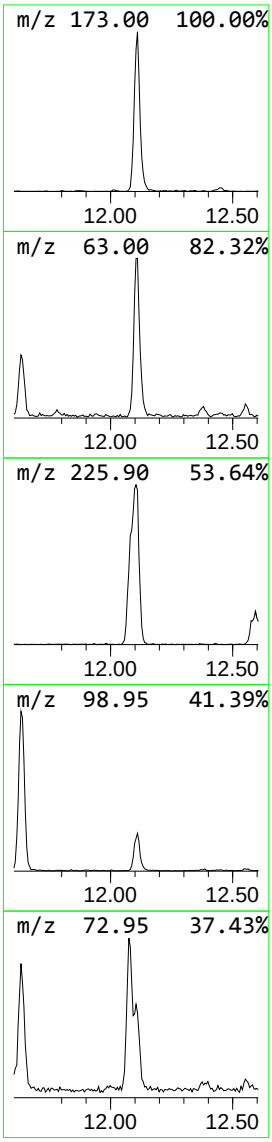
****

Peak Number 10 4-Chloro-2-nitrophenol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.104	2.54 ng/ul	506905	Naphthalene-d8	10.834

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	99
2	4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	96
3	4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	96
4	2-Chloro-6-nitrophenol	173	C6H4ClNO3	000603-86-1	91
5	Phenol, 5-chloro-2-nitro-	173	C6H4ClNO3	000611-07-4	81





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120222\
Data File : BP012942.D
Acq On : 02 Dec 2022 16:35
Operator : CG/JU
Sample : N5785-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L75

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

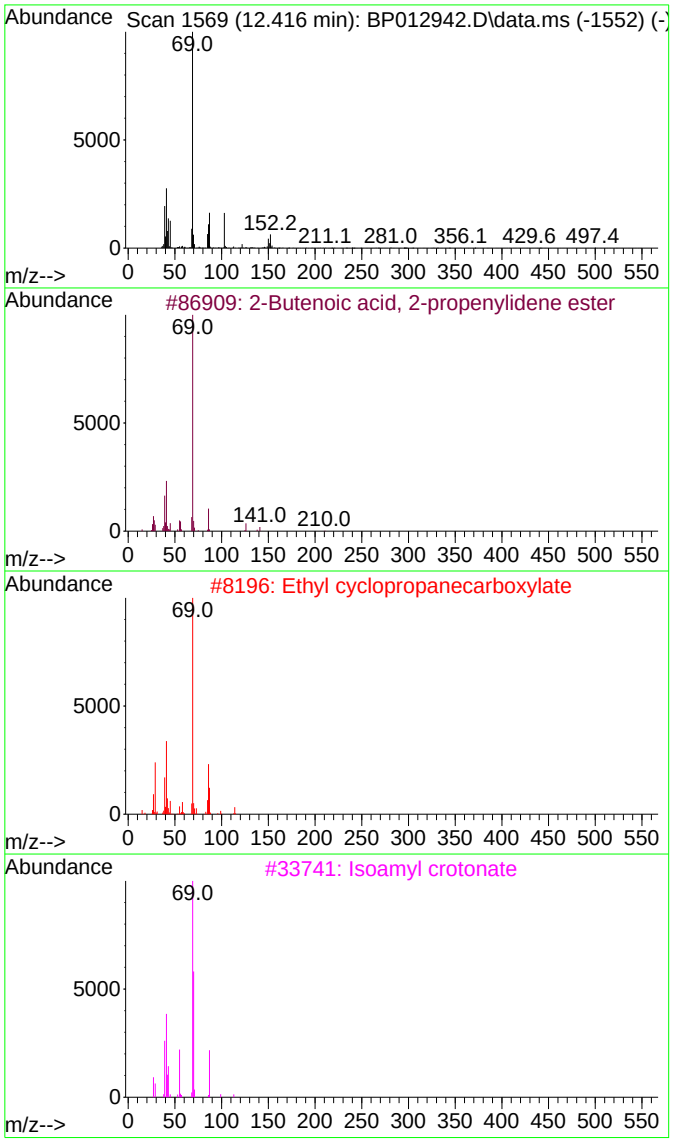
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

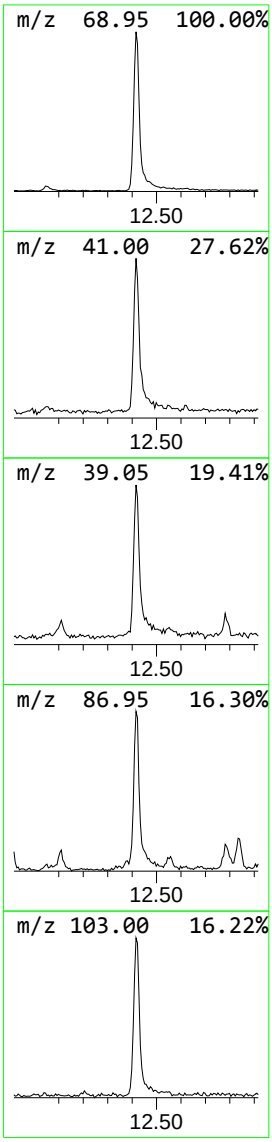
****

Peak Number 11 unknown-01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.416	2.56 ng/ul	510531	Naphthalene-d8	10.834

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butenoic acid, 2-propenylidene...	210	C11H14O4	055030-70-1	40
2	Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	38
3	Isoamyl crotonate	156	C9H16O2	1010429-17-3	33
4	Cyclopropanecarboxylic acid, 3-m...	154	C9H14O2	1000299-37-4	33
5	Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120222\
Data File : BP012942.D
Acq On : 02 Dec 2022 16:35
Operator : CG/JU
Sample : N5785-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L75

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

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

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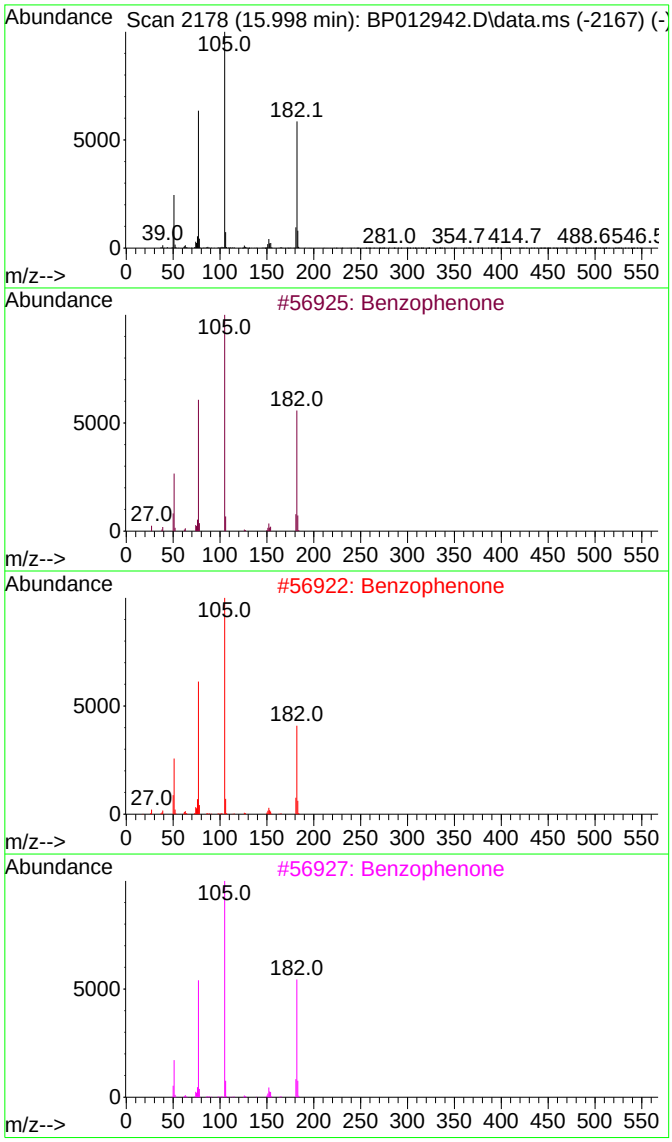
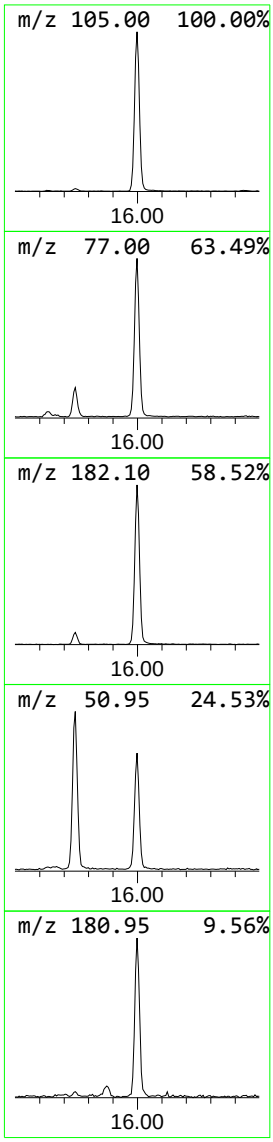
Peak Number 12 Benzophenone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.998	2.61 ng/ul	851494	Acenaphthene-d10	14.639

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	Benzophenone	182	C13H10O	000119-61-9	95
3	Benzophenone	182	C13H10O	000119-61-9	95
4	Benzophenone	182	C13H10O	000119-61-9	93
5	Benzophenone	182	C13H10O	000119-61-9	91





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Data File : BP012942.D
Acq On : 02 Dec 2022 16:35
Operator : CG/JU
Sample : N5785-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L75

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

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

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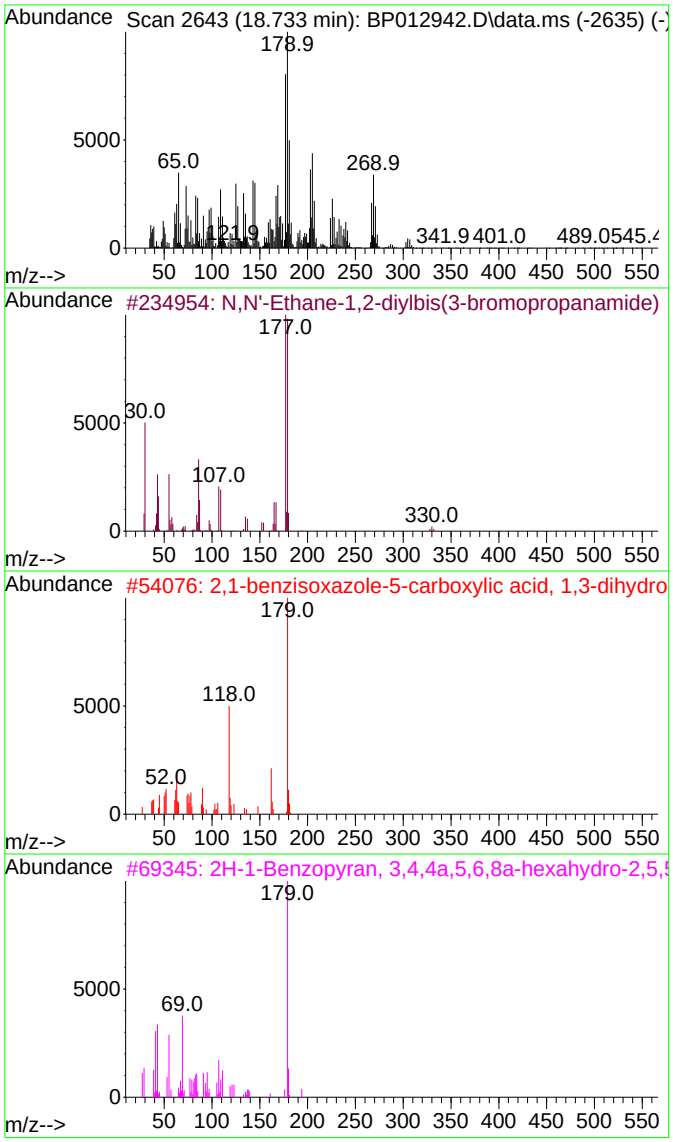
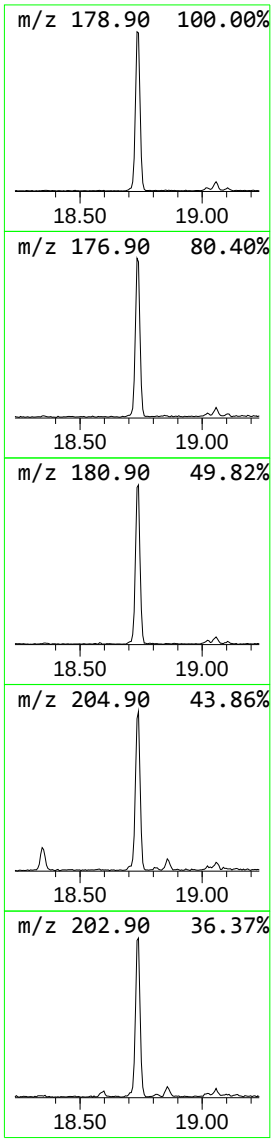
Peak Number 13 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.733	5.86 ng/ul	2572800	Phenanthrene-d10	17.398

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N,N'-Ethane-1,2-diylbis(3-bromop...	328	C8H14Br2N2O2	016044-35-2	35
2	2,1-benzisoxazole-5-carboxylic a...	179	C8H5NO4	1000404-80-3	30
3	2H-1-Benzopyran, 3,4,4a,5,6,8a-h...	194	C13H22O	041678-32-4	25
4	Phenol, 4-amino-2,6-dichloro-	177	C6H5Cl2NO	005930-28-9	25
5	6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	25





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Data File : BP012942.D
Acq On : 02 Dec 2022 16:35
Operator : CG/JU
Sample : N5785-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L75

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

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

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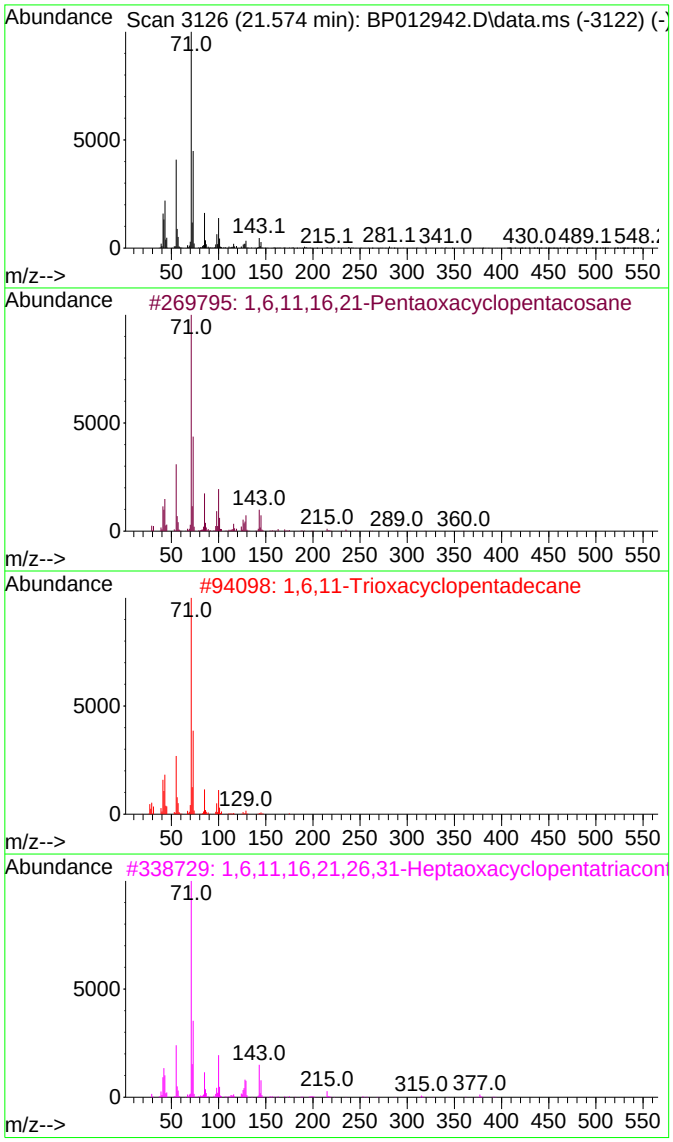
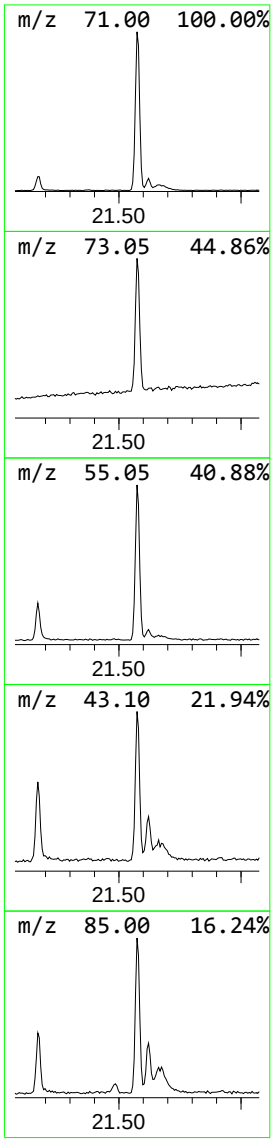
Peak Number 14 1,6,11,16,21-Pentaoxacyclop... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.574	2.24 ng/ul	1270590	Chrysene-d12	21.480

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,6,11,16,21-Pentaoxacyclopentac...	360	C20H40O5	056890-57-4	86
2	1,6,11-Trioxacyclopentadecane	216	C12H24O3	000295-63-6	86
3	1,6,11,16,21,26,31-Heptaoxacyclo...	504	C28H56O7	066055-34-3	83
4	1,6,11,16-Tetraoxacycloeicosane	288	C16H32O4	017043-02-6	78
5	1,6,11,16,21,26-Hexaoxacyclotria...	432	C24H48O6	064001-05-4	62





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Acq On : 02 Dec 2022 16:35
Operator : CG/JU
Sample : N5785-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L75

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

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

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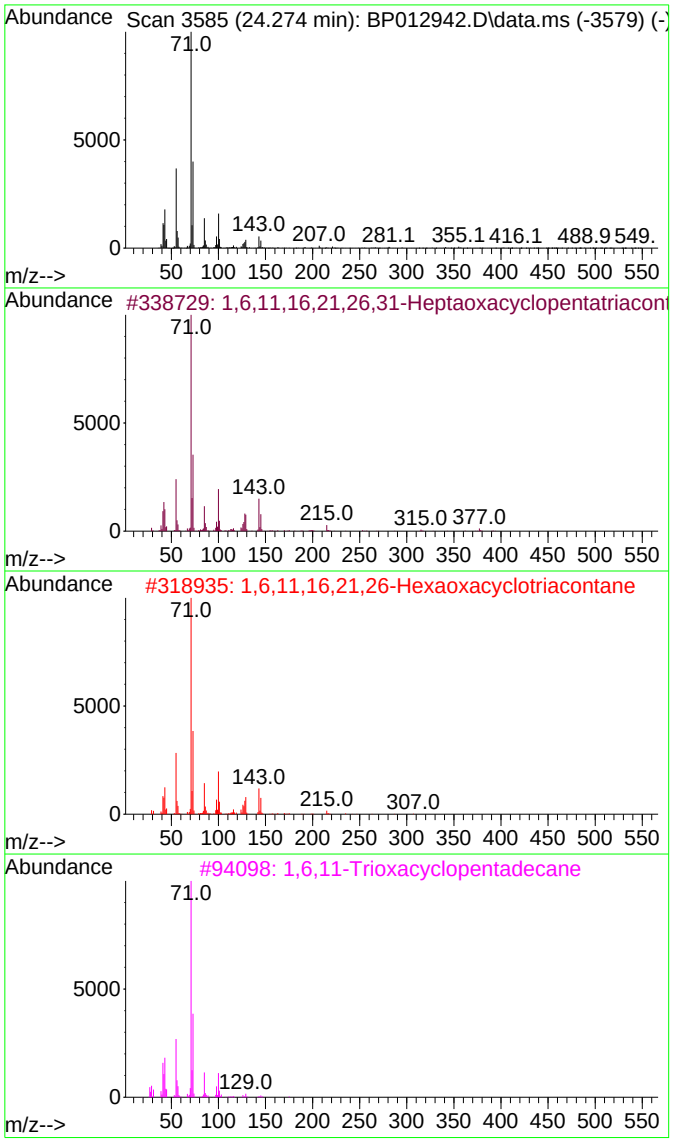
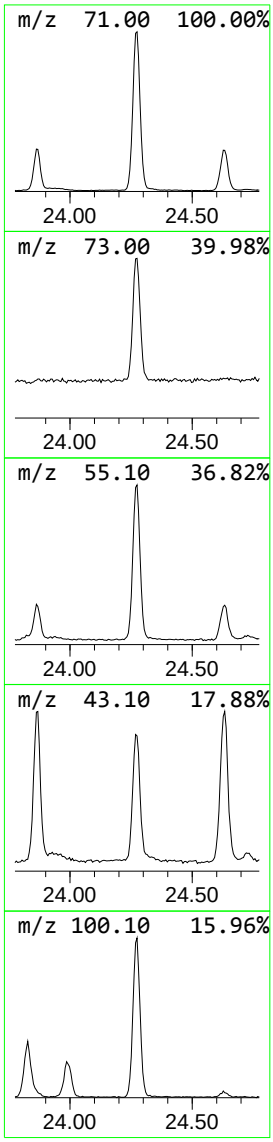
Peak Number 15 1,6,11,16,21,26,31-Heptaoxa... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.274	3.78 ng/ul	2127130	Perylene-d12	23.992

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,6,11,16,21,26,31-Heptaoxacyclo...	504	C28H56O7	066055-34-3	91
2	1,6,11,16,21,26-Hexaoxacyclotria...	432	C24H48O6	064001-05-4	91
3	1,6,11-Trioxacyclopentadecane	216	C12H24O3	000295-63-6	86
4	1,6,11,16,21-Pentaoxacyclopentac...	360	C20H40O5	056890-57-4	52
5	Acrolein,dimethyl acetal	102	C5H10O2	006044-68-4	50





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Operator : CG/JU
Sample : N5785-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L75

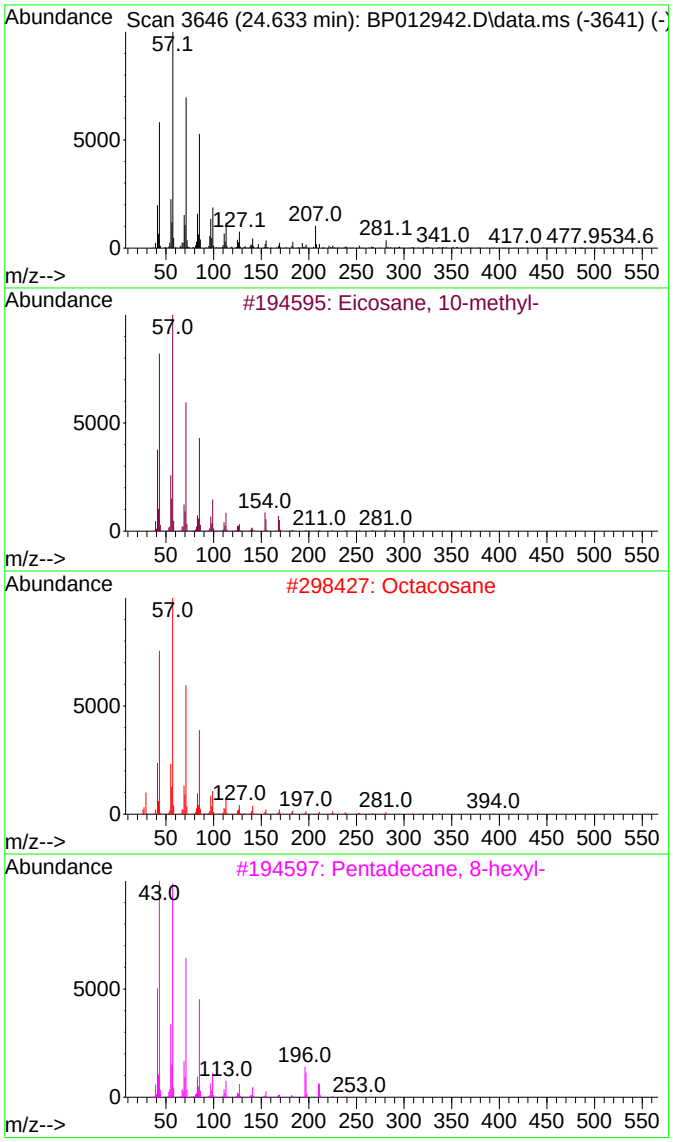
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

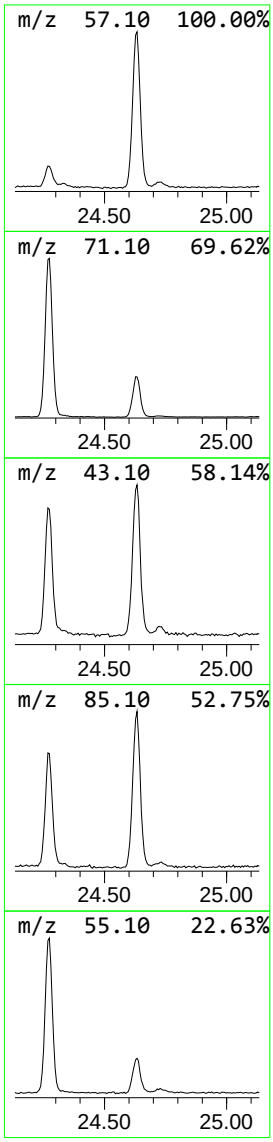
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.633	2.15 ng/ul	1209010	Perylene-d12	23.992

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 10-methyl-	296	C21H44	054833-23-7	91
2	Octacosane	394	C28H58	000630-02-4	87
3	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	87
4	Heneicosane	296	C21H44	000629-94-7	87
5	Octacosane, 2-methyl-	408	C29H60	001560-98-1	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120222\
Data File : BP012942.D
Acq On : 02 Dec 2022 16:35
Operator : CG/JU
Sample : N5785-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0L75

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.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, 1-brom...	9.510	7.8	ng/ul	1559350	2	10.834	3995460	20.0
Benzene, 1-brom...	9.828	4.8	ng/ul	954417	2	10.834	3995460	20.0
Benzene, 1-chlo...	11.422	17.7	ng/ul	3541570	2	10.834	3995460	20.0
Benzene, 1-chlo...	11.634	4.5	ng/ul	892018	2	10.834	3995460	20.0
4-Chloro-2-nitr...	12.104	2.5	ng/ul	506905	2	10.834	3995460	20.0
unknown-01	12.416	2.6	ng/ul	510531	2	10.834	3995460	20.0
Benzophenone	15.998	2.6	ng/ul	851494	3	14.639	6536390	20.0
unknown-02	18.733	5.9	ng/ul	2572800	4	17.398	8785270	20.0
1,6,11,16,21-Pe...	21.574	2.2	ng/ul	1270590	5	21.480	11369300	20.0
1,6,11,16,21,26...	24.274	3.8	ng/ul	2127130	6	23.992	11254200	20.0
(DEL) Alkane: S...	24.633	2.1	ng/ul	1209010	6	23.992	11254200	20.0