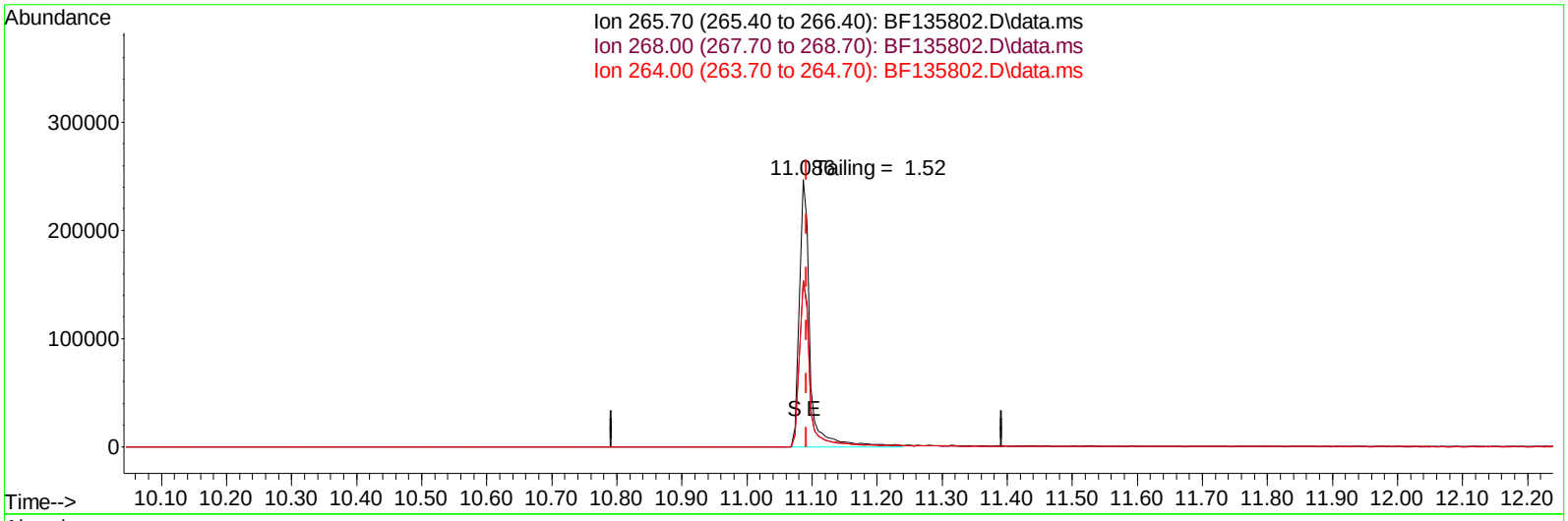
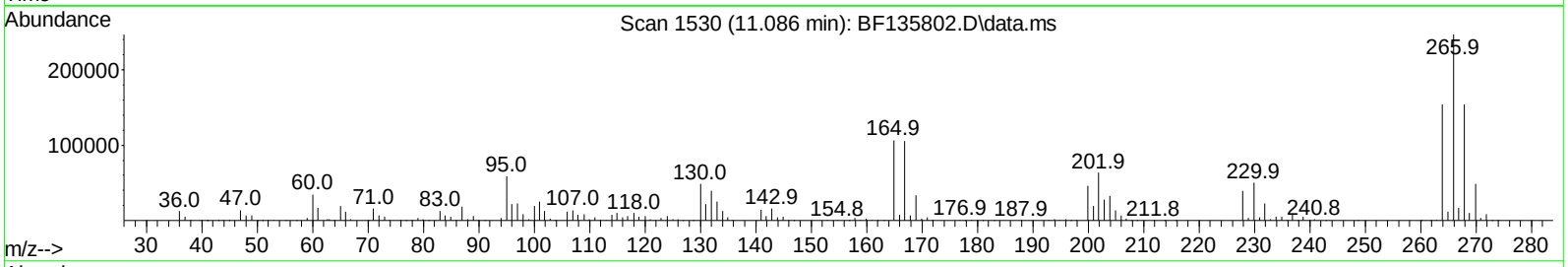
Quantitation Report (Qedit)

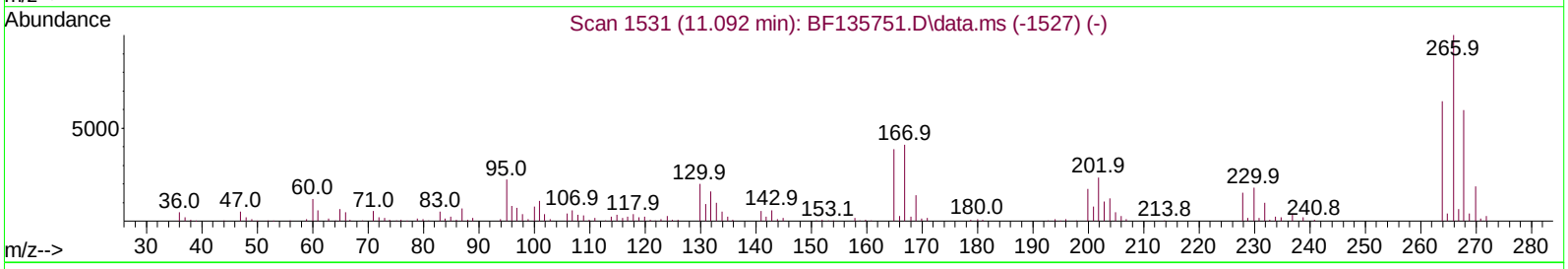
Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101723\
Data File : BF135802.D
Acq On : 17 Oct 2023 09:54
Operator : CG\JU
Sample : DFTPP
Misc :
ALS Vial : 1 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SSTD0.4299

Quant Time: Oct 18 05:23:51 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 17 01:22:00 2023
Response via : Initial Calibration







TIC: BF135802.D\data.ms

(70) Pentachlorophenol (C)

11.086min (-0.006) 325830.29 ng

response 271868

Ion	Exp%	Act%
265.70	100.00	100.00
268.00	59.60	62.40
264.00	64.30	62.34
0.00	0.00	0.00

Quantitation Report (Not Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101823\
Data File : BN028316.D
Acq On : 18 Oct 2023 09:49
Operator : MA/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SSTD0.4299

Quant Time: Oct 18 10:59:59 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN100323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 16 16:16:04 2023
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.882	152	2611	0.400	ng/ul	0.00
4) Naphthalene-d8	10.708	136	8451	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.532	164	4970	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.299	188	11158	0.400	ng/ul	0.00
17) Chrysene-d12	21.492	240	11624	0.400	ng/ul	0.00
23) Perylene-d12	23.942	264	12785	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.263	96	1370	0.418	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.297	152	5074	0.373	ng/ul	0.00
18) Fluoranthene-d10	19.324	212	12869	0.349	ng/ul	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.296	88	1420	0.409	ng/ul#	91
5) Naphthalene	10.757	128	9171	0.382	ng/ul	99
7) 2-Methylnaphthalene	12.368	142	5666	0.374	ng/ul	99
8) 1-Methylnaphthalene	12.583	142	5808	0.371	ng/ul	99
10) Acenaphthylene	14.259	152	8206	0.364	ng/ul	100
11) Acenaphthene	14.597	153	6865	0.383	ng/ul	99
12) Fluorene	15.591	166	7901	0.396	ng/ul	98
14) Pentachlorophenol	16.949	266	691	0.289	ng/ul	96
15) Phenanthrene	17.341	178	13049	0.392	ng/ul	99
16) Anthracene	17.434	178	11352	0.382	ng/ul	99
19) Fluoranthene	19.357	202	15643	0.352	ng/ul	100
20) Pyrene	19.724	202	16305	0.354	ng/ul	99
21) Benzo(a)anthracene	21.475	228	16035	0.378	ng/ul	100
22) Chrysene	21.530	228	16629	0.388	ng/ul	100
24) Benzo(b)fluoranthene	23.181	252	18556	0.391	ng/ul	98
25) Benzo(k)fluoranthene	23.231	252	18574	0.394	ng/ul	99
26) Benzo(a)pyrene	23.831	252	15806	0.391	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.508	276	21320	0.436	ng/ul	99
28) Dibenzo(a,h)anthracene	26.518	278	17241	0.433	ng/ul	98
29) Benzo(g,h,i)perylene	27.317	276	18101	0.453	ng/ul	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101823\
Data File : BN028316.D
Acq On : 18 Oct 2023 09:49
Operator : MA/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SSTD0.4299

Quant Time: Oct 18 10:59:59 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN100323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 16 16:16:04 2023
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