

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121624\
 Data File : BN035616.D
 Acq On : 16 Dec 2024 14:37
 Operator : RC/JU
 Sample : SST0.846
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.8233

Quant Time: Dec 16 15:16:13 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN121624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 16 14:42:43 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.288	152	1893	0.400	ng/ul	0.00
4) Naphthalene-d8	10.031	136	5607	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.949	164	3947	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.717	188	8816	0.400	ng/ul	0.00
17) Chrysene-d12	20.964	240	7780	0.400	ng/ul	0.00
23) Perylene-d12	23.054	264	8450	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.955	96	1355	0.791	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.637	152	6285	0.775	ng/ul	0.00
18) Fluoranthene-d10	18.767	212	17830	0.831	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	2.984	88	1381	0.741	ng/ul#	85
5) Naphthalene	10.080	128	10998	0.775	ng/ul	97
7) 2-Methylnaphthalene	11.714	142	7295	0.759	ng/ul	99
8) 1-Methylnaphthalene	11.939	142	7517	0.770	ng/ul	100
10) Acenaphthylene	13.657	152	12072	0.775	ng/ul	98
11) Acenaphthene	14.009	153	9075	0.763	ng/ul	100
12) Fluorene	15.013	166	10607	0.761	ng/ul	100
14) Pentachlorophenol	16.384	266	1696	0.809	ng/ul	99
15) Phenanthrene	16.759	178	17533	0.763	ng/ul	98
16) Anthracene	16.848	178	16389	0.766	ng/ul	99
19) Fluoranthene	18.800	202	21788	0.810	ng/ul	98
20) Pyrene	19.167	202	23507	0.830	ng/ul	98
21) Benzo(a)anthracene	20.947	228	20314	0.763	ng/ul	99
22) Chrysene	20.999	228	20824	0.771	ng/ul	99
24) Benzo(b)fluoranthene	22.440	252	19742	0.724	ng/ul	92
25) Benzo(k)fluoranthene	22.480	252	21801	0.770	ng/ul#	92
26) Benzo(a)pyrene	22.963	252	19028	0.757	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	25.081	276	23595	0.709	ng/ul#	85
28) Dibenzo(a,h)anthracene	25.098	278	18221	0.702	ng/ul#	92
29) Benzo(g,h,i)perylene	25.691	276	18643	0.709	ng/ul	96

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

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