

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121621\
 Data File : BN018025.D
 Acq On : 15 Dec 2021 15:15
 Operator : CG/JU
 Sample : SST0.445
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.4245

Quant Time: Dec 15 17:20:14 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN121621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 15 17:18:54 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.666	152	4803	0.400	ng/ul	0.00
4) Naphthalene-d8	10.465	136	13928	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.314	164	8896	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.067	188	17795	0.400	ng/ul #	0.00
17) Chrysene-d12	21.265	240	18750	0.400	ng/ul #	0.00
23) Perylene-d12	23.535	264	15824	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.143	96	2011	0.382	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.066	152	6420	0.299	ng/ul	0.00
18) Fluoranthene-d10	19.102	212	21402	0.318	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.177	88	2378	0.350	ng/ul#	45
5) Naphthalene	10.509	128	14613	0.379	ng/ul	98
7) 2-Methylnaphthalene	12.137	142	8396	0.304	ng/ul	100
8) 1-Methylnaphthalene	12.352	142	9673	0.355	ng/ul	100
10) Acenaphthylene	14.032	152	13749	0.322	ng/ul	99
11) Acenaphthene	14.374	153	12347	0.404	ng/ul	97
12) Fluorene	15.374	166	13718	0.379	ng/ul	99
14) Pentachlorophenol	16.755	266	996	0.256	ng/ul	97
15) Phenanthrene	17.109	178	23390	0.437	ng/ul	99
16) Anthracene	17.206	178	18566	0.345	ng/ul	99
19) Fluoranthene	19.129	202	30139	0.347	ng/ul	98
20) Pyrene	19.492	202	33334	0.375	ng/ul#	94
21) Benzo(a)anthracene	21.250	228	20760	0.298	ng/ul	99
22) Chrysene	21.300	228	31504	0.451	ng/ul	98
24) Benzo(b)fluoranthene	22.851	252	25402	0.393	ng/ul	85
25) Benzo(k)fluoranthene	22.892	252	31459	0.472	ng/ul#	83
26) Benzo(a)pyrene	23.439	252	23241	0.430	ng/ul#	81
27) Indeno(1,2,3-cd)pyrene	25.875	276	25474	0.559	ng/ul#	90
28) Dibenzo(a,h)anthracene	25.898	278	18550	0.547	ng/ul#	83
29) Benzo(g,h,i)perylene	26.576	276	23037	0.553	ng/ul#	89

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

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