

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020222\
 Data File : BN018543.D
 Acq On : 02 Feb 2022 11:11
 Operator : CG/JU
 Sample : SSTD0.237
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.2237

Quant Time: Feb 02 13:44:46 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN020222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 02 13:41:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.927	152	486	0.400	ng/ul	0.00
4) Naphthalene-d8	10.733	136	1635	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.558	164	872	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.294	188	1653	0.400	ng/ul #	0.00
17) Chrysene-d12	21.465	240	1564	0.400	ng/ul #	0.00
23) Perylene-d12	23.859	264	1422	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	120	0.214	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.323	152	447	0.203	ng/ul	0.00
18) Fluoranthene-d10	19.314	212	913	0.209	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.337	88	124	0.211	ng/ul#	88
5) Naphthalene	10.783	128	918	0.204	ng/ul#	85
7) 2-Methylnaphthalene	12.394	142	578	0.202	ng/ul	99
8) 1-Methylnaphthalene	12.609	142	601	0.206	ng/ul	100
10) Acenaphthylene	14.276	152	697	0.192	ng/ul#	81
11) Acenaphthene	14.618	153	621	0.206	ng/ul	97
12) Fluorene	15.604	166	652	0.202	ng/ul#	96
14) Pentachlorophenol	16.947	266	73	0.497	ng/ul	97
15) Phenanthrene	17.336	178	1087	0.204	ng/ul#	88
16) Anthracene	17.429	178	849	0.193	ng/ul#	84
19) Fluoranthene	19.346	202	1238	0.205	ng/ul#	80
20) Pyrene	19.704	202	1305	0.210	ng/ul#	80
21) Benzo(a)anthracene	21.447	228	1000	0.194	ng/ul#	91
22) Chrysene	21.500	228	1228	0.207	ng/ul	93
24) Benzo(b)fluoranthene	23.128	252	1216	0.205	ng/ul#	82
25) Benzo(k)fluoranthene	23.174	252	1302	0.209	ng/ul#	81
26) Benzo(a)pyrene	23.753	252	1006	0.204	ng/ul#	71
27) Indeno(1,2,3-cd)pyrene	26.346	276	1313	0.204	ng/ul#	81
28) Dibenzo(a,h)anthracene	26.363	278	1016	0.201	ng/ul#	69
29) Benzo(g,h,i)perylene	27.101	276	1151	0.206	ng/ul#	71

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

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