

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020222\
 Data File : BN018546.D
 Acq On : 02 Feb 2022 12:58
 Operator : CG/JU
 Sample : SSTD1.640
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.6240

Quant Time: Feb 02 13:45:20 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	413	0.400	ng/ul	0.00
4) Naphthalene-d8	10.733	136	1447	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.558	164	762	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.294	188	1482	0.400	ng/ul #	0.00
17) Chrysene-d12	21.462	240	1389	0.400	ng/ul #	0.00
23) Perylene-d12	23.858	264	1116	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.294	96	708	1.488	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.317	152	3029	1.552	ng/ul	0.00
18) Fluoranthene-d10	19.314	212	5845	1.506	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.332	88	770	1.540	ng/ul#	72
5) Naphthalene	10.783	128	6027	1.513	ng/ul	100
7) 2-Methylnaphthalene	12.389	142	3981	1.574	ng/ul	100
8) 1-Methylnaphthalene	12.609	142	3984	1.540	ng/ul	100
10) Acenaphthylene	14.276	152	4603	1.453	ng/ul	98
11) Acenaphthene	14.618	153	3992	1.516	ng/ul	99
12) Fluorene	15.599	166	4534	1.608	ng/ul	99
14) Pentachlorophenol	16.943	266	382	2.901	ng/ul	99
15) Phenanthrene	17.336	178	7310	1.530	ng/ul	99
16) Anthracene	17.424	178	6239	1.582	ng/ul	99
19) Fluoranthene	19.346	202	8364	1.559	ng/ul#	93
20) Pyrene	19.704	202	8289	1.504	ng/ul#	90
21) Benzo(a)anthracene	21.447	228	6826	1.493	ng/ul	99
22) Chrysene	21.500	228	7999	1.518	ng/ul	99
24) Benzo(b)fluoranthene	23.128	252	7529	1.618	ng/ul	97
25) Benzo(k)fluoranthene	23.174	252	8016	1.643	ng/ul#	96
26) Benzo(a)pyrene	23.753	252	6191	1.599	ng/ul#	96
27) Indeno(1,2,3-cd)pyrene	26.343	276	8251	1.636	ng/ul#	81
28) Dibenzo(a,h)anthracene	26.366	278	6569	1.658	ng/ul#	93
29) Benzo(g,h,i)perylene	27.104	276	7182	1.636	ng/ul#	90

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

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