

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN010422\
 Data File : BN018330.D
 Acq On : 04 Jan 2022 11:19
 Operator : CG/JU
 Sample : SSTD0.417
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4217

Quant Time: Jan 04 12:34:30 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN010422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 04 12:28:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.603	152	5498	0.400	ng/ul	0.00
4) Naphthalene-d8	10.383	136	18136	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.240	164	12221	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.987	188	29029	0.400	ng/ul	0.00
17) Chrysene-d12	21.183	240	32871	0.400	ng/ul	0.00
23) Perylene-d12	23.398	264	29832	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.067	96	2064	0.402	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.983	152	9715	0.308	ng/ul	0.00
18) Fluoranthene-d10	19.018	212	34470	0.354	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.101	88	2315	0.405	ng/ul#	65
5) Naphthalene	10.432	128	19405	0.361	ng/ul	99
7) 2-Methylnaphthalene	12.060	142	12490	0.317	ng/ul	99
8) 1-Methylnaphthalene	12.275	142	13254	0.334	ng/ul	99
10) Acenaphthylene	13.963	152	17773	0.337	ng/ul	99
11) Acenaphthene	14.305	153	16196	0.381	ng/ul	100
12) Fluorene	15.295	166	19390	0.369	ng/ul	99
14) Pentachlorophenol	16.658	266	2476	0.215	ng/ul	99
15) Phenanthrene	17.029	178	35581	0.371	ng/ul	99
16) Anthracene	17.122	178	28661	0.332	ng/ul	99
19) Fluoranthene	19.051	202	46520	0.356	ng/ul	100
20) Pyrene	19.413	202	48652	0.356	ng/ul	99
21) Benzo(a)anthracene	21.165	228	38867	0.310	ng/ul	100
22) Chrysene	21.218	228	50720	0.373	ng/ul	100
24) Benzo(b)fluoranthene	22.734	252	46365	0.371	ng/ul	100
25) Benzo(k)fluoranthene	22.778	252	54651	0.384	ng/ul	100
26) Benzo(a)pyrene	23.305	252	41509	0.351	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.654	276	53828	0.357	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.674	278	40899	0.351	ng/ul	99
29) Benzo(g,h,i)perylene	26.335	276	49212	0.384	ng/ul	99

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

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