

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102122\
 Data File : BN022256.D
 Acq On : 21 Oct 2022 16:17
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4321

Quant Time: Oct 21 22:40:11 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN102022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 19 23:32:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.949	152	4130	0.400	ng/ul	0.00
4) Naphthalene-d8	10.778	136	13699	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.623	164	7699	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.379	188	14565	0.400	ng/ul	0.00
17) Chrysene-d12	21.579	240	8543	0.400	ng/ul	0.00
23) Perylene-d12	24.040	264	6783	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.257	96	1931	0.360	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.373	152	8176	0.394	ng/ul	0.00
18) Fluoranthene-d10	19.412	212	13559	0.473	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.295	88	1967	0.357	ng/ul	95
5) Naphthalene	10.827	128	15818	0.391	ng/ul	100
7) 2-Methylnaphthalene	12.444	142	10116	0.397	ng/ul	99
8) 1-Methylnaphthalene	12.664	142	10399	0.396	ng/ul	100
10) Acenaphthylene	14.341	152	14099	0.408	ng/ul	99
11) Acenaphthene	14.683	153	11206	0.384	ng/ul	100
12) Fluorene	15.674	166	12223	0.361	ng/ul	99
14) Pentachlorophenol	17.032	266	1129	0.326	ng/ul	100
15) Phenanthrene	17.421	178	18774	0.384	ng/ul	100
16) Anthracene	17.514	178	16260	0.395	ng/ul	100
19) Fluoranthene	19.445	202	18259	0.467	ng/ul	98
20) Pyrene	19.807	202	18326	0.466	ng/ul	100
21) Benzo(a)anthracene	21.562	228	13246	0.414	ng/ul	99
22) Chrysene	21.617	228	13165	0.388	ng/ul	99
24) Benzo(b)fluoranthene	23.280	252	11941	0.351	ng/ul	94
25) Benzo(k)fluoranthene	23.330	252	11759	0.355	ng/ul	95
26) Benzo(a)pyrene	23.926	252	10340	0.381	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	26.622	276	13152	0.390	ng/ul	100
28) Dibenzo(a,h)anthracene	26.645	278	10341	0.383	ng/ul	97
29) Benzo(g,h,i)perylene	27.417	276	10831	0.394	ng/ul	99

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

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