

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111924\
 Data File : BN035181.D
 Acq On : 20 Nov 2024 03:16
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4344

Quant Time: Nov 20 03:48:03 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN111624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 18 00:44:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.541	152	3323	0.400	ng/ul	0.00
4) Naphthalene-d8	10.306	136	7986	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.185	164	5570	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.937	188	11597	0.400	ng/ul	0.00
17) Chrysene-d12	21.145	240	8751	0.400	ng/ul	0.00
23) Perylene-d12	23.320	264	9394	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.136	96	1272	0.419	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.906	152	4931	0.421	ng/ul	0.00
18) Fluoranthene-d10	18.977	212	10623	0.440	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.170	88	1354	0.411	ng/ul	98
5) Naphthalene	10.355	128	8363	0.416	ng/ul	99
7) 2-Methylnaphthalene	11.978	142	5935	0.425	ng/ul	99
8) 1-Methylnaphthalene	12.203	142	6009	0.427	ng/ul	99
10) Acenaphthylene	13.898	152	8874	0.402	ng/ul	99
11) Acenaphthene	14.245	153	6897	0.408	ng/ul	98
12) Fluorene	15.240	166	8086	0.408	ng/ul	99
14) Pentachlorophenol	16.595	266	886	0.328	ng/ul	98
15) Phenanthrene	16.979	178	12449	0.406	ng/ul	100
16) Anthracene	17.068	178	11346	0.399	ng/ul	100
19) Fluoranthene	19.005	202	13445	0.441	ng/ul	100
20) Pyrene	19.372	202	13806	0.435	ng/ul	100
21) Benzo(a)anthracene	21.131	228	12198	0.402	ng/ul	100
22) Chrysene	21.183	228	12656	0.414	ng/ul	99
24) Benzo(b)fluoranthene	22.671	252	12596	0.406	ng/ul	99
25) Benzo(k)fluoranthene	22.714	252	13260	0.414	ng/ul	98
26) Benzo(a)pyrene	23.223	252	11433	0.403	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.473	276	15746	0.411	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.490	278	12274	0.409	ng/ul	98
29) Benzo(g,h,i)perylene	26.127	276	12275	0.406	ng/ul	99

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

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