

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111924\
 Data File : BN035157.D
 Acq On : 19 Nov 2024 12:05
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4342

Quant Time: Nov 19 12:32:32 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	7.545	152	2627	0.400	ng/ul	0.00
4) Naphthalene-d8	10.306	136	6138	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.185	164	4148	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.941	188	8727	0.400	ng/ul	0.00
17) Chrysene-d12	21.151	240	7280	0.400	ng/ul	0.00
23) Perylene-d12	23.325	264	7905	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	1020	0.425	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.906	152	3738	0.416	ng/ul	0.00
18) Fluoranthene-d10	18.977	212	8292	0.413	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.174	88	1048	0.402	ng/ul	96
5) Naphthalene	10.355	128	6433	0.416	ng/ul	99
7) 2-Methylnaphthalene	11.978	142	4523	0.422	ng/ul	100
8) 1-Methylnaphthalene	12.203	142	4544	0.420	ng/ul	99
10) Acenaphthylene	13.902	152	6747	0.410	ng/ul	99
11) Acenaphthene	14.249	153	5188	0.412	ng/ul	99
12) Fluorene	15.244	166	5986	0.405	ng/ul	99
14) Pentachlorophenol	16.595	266	690	0.339	ng/ul	97
15) Phenanthrene	16.979	178	9422	0.408	ng/ul	99
16) Anthracene	17.072	178	8559	0.400	ng/ul	100
19) Fluoranthene	19.009	202	10690	0.422	ng/ul	99
20) Pyrene	19.372	202	10909	0.414	ng/ul	98
21) Benzo(a)anthracene	21.136	228	10174	0.403	ng/ul	100
22) Chrysene	21.186	228	10584	0.416	ng/ul	100
24) Benzo(b)fluoranthene	22.679	252	10814	0.414	ng/ul	99
25) Benzo(k)fluoranthene	22.720	252	11415	0.424	ng/ul	99
26) Benzo(a)pyrene	23.229	252	9760	0.409	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.483	276	13541	0.420	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.500	278	10535	0.417	ng/ul	99
29) Benzo(g,h,i)perylene	26.137	276	10685	0.420	ng/ul	100

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

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