

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111824\
 Data File : BN035142.D
 Acq On : 18 Nov 2024 10:57
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4340

Quant Time: Nov 18 11:47:39 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	2671	0.400	ng/ul	0.00
4) Naphthalene-d8	10.311	136	6130	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.189	164	4050	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.941	188	8730	0.400	ng/ul	0.00
17) Chrysene-d12	21.151	240	7863	0.400	ng/ul	0.00
23) Perylene-d12	23.323	264	8507	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	999	0.409	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.906	152	3672	0.409	ng/ul	0.00
18) Fluoranthene-d10	18.981	212	8607	0.397	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.174	88	1092	0.412	ng/ul	93
5) Naphthalene	10.361	128	6412	0.415	ng/ul	100
7) 2-Methylnaphthalene	11.983	142	4394	0.410	ng/ul	98
8) 1-Methylnaphthalene	12.203	142	4434	0.410	ng/ul	98
10) Acenaphthylene	13.902	152	6552	0.408	ng/ul	100
11) Acenaphthene	14.249	153	5010	0.408	ng/ul	99
12) Fluorene	15.244	166	5800	0.402	ng/ul	100
14) Pentachlorophenol	16.599	266	725	0.357	ng/ul	98
15) Phenanthrene	16.983	178	9377	0.406	ng/ul	99
16) Anthracene	17.072	178	8523	0.398	ng/ul	99
19) Fluoranthene	19.009	202	10924	0.399	ng/ul	100
20) Pyrene	19.372	202	11343	0.398	ng/ul	97
21) Benzo(a)anthracene	21.134	228	10898	0.399	ng/ul	100
22) Chrysene	21.186	228	11246	0.410	ng/ul	99
24) Benzo(b)fluoranthene	22.676	252	11444	0.407	ng/ul	100
25) Benzo(k)fluoranthene	22.717	252	11945	0.412	ng/ul	99
26) Benzo(a)pyrene	23.229	252	10410	0.405	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.480	276	14106	0.406	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.493	278	10889	0.400	ng/ul	99
29) Benzo(g,h,i)perylene	26.134	276	11167	0.408	ng/ul	99

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

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