

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071224\
 Data File : BN032636.D
 Acq On : 11 Jul 2024 15:18
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4254

Quant Time: Jul 12 01:54:51 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN071224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 01:40:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.591	152	3593	0.400	ng/ul	0.01
4) Naphthalene-d8	10.365	136	12265	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.221	164	7904	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.991	188	17452	0.400	ng/ul	0.00
17) Chrysene-d12	21.200	240	13981	0.400	ng/ul	0.00
23) Perylene-d12	23.403	264	14094	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.106	96	1910	0.443	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.977	152	7257	0.391	ng/ul	0.00
18) Fluoranthene-d10	19.027	212	17879	0.392	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.139	88	2260	0.460	ng/ul	98
5) Naphthalene	10.415	128	12916	0.392	ng/ul	99
7) 2-Methylnaphthalene	12.054	142	8228	0.400	ng/ul	99
8) 1-Methylnaphthalene	12.252	142	8512	0.400	ng/ul	98
10) Acenaphthylene	13.943	152	12550	0.393	ng/ul	99
11) Acenaphthene	14.281	153	10436	0.395	ng/ul	99
12) Fluorene	15.299	166	11975	0.397	ng/ul	99
14) Pentachlorophenol	16.678	266	1042	0.340	ng/ul	99
15) Phenanthrene	17.033	178	18548	0.390	ng/ul	100
16) Anthracene	17.143	178	14968	0.385	ng/ul	100
19) Fluoranthene	19.059	202	22315	0.393	ng/ul	99
20) Pyrene	19.426	202	23040	0.392	ng/ul	99
21) Benzo(a)anthracene	21.188	228	16904	0.380	ng/ul	99
22) Chrysene	21.238	228	25248	0.392	ng/ul	100
24) Benzo(b)fluoranthene	22.740	252	18655	0.375	ng/ul	100
25) Benzo(k)fluoranthene	22.784	252	23714	0.391	ng/ul	99
26) Benzo(a)pyrene	23.313	252	16269	0.378	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.633	276	21711	0.373	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.646	278	16640	0.368	ng/ul	99
29) Benzo(g,h,i)perylene	26.307	276	18130	0.372	ng/ul	100

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

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