

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090324\
 Data File : BN033734.D
 Acq On : 03 Sep 2024 10:55
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4302

Quant Time: Sep 04 01:37:17 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN082724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 04 01:36:40 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.523	152	6679	0.400	ng/ul	0.00
4) Naphthalene-d8	10.283	136	17645	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.161	164	8595	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.915	188	17489	0.400	ng/ul	0.00
17) Chrysene-d12	21.118	240	14313	0.400	ng/ul	0.00
23) Perylene-d12	23.275	264	15621	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.135	96	2870	0.418	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.878	152	9753	0.385	ng/ul	0.00
18) Fluoranthene-d10	18.952	212	16983	0.398	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.169	88	3132	0.411	ng/ul#	87
5) Naphthalene	10.332	128	18457	0.391	ng/ul	99
7) 2-Methylnaphthalene	11.955	142	11429	0.382	ng/ul	99
8) 1-Methylnaphthalene	12.175	142	11803	0.384	ng/ul	99
10) Acenaphthylene	13.874	152	15020	0.385	ng/ul	100
11) Acenaphthene	14.225	153	11442	0.406	ng/ul	99
12) Fluorene	15.215	166	12669	0.402	ng/ul	99
14) Pentachlorophenol	16.568	266	1778	0.323	ng/ul	99
15) Phenanthrene	16.953	178	19616	0.380	ng/ul	99
16) Anthracene	17.046	178	17252	0.372	ng/ul	100
19) Fluoranthene	18.980	202	21766	0.385	ng/ul	99
20) Pyrene	19.343	202	23123	0.386	ng/ul	99
21) Benzo(a)anthracene	21.103	228	21107	0.375	ng/ul	99
22) Chrysene	21.156	228	21587	0.379	ng/ul	100
24) Benzo(b)fluoranthene	22.634	252	22456	0.370	ng/ul	98
25) Benzo(k)fluoranthene	22.678	252	23551	0.388	ng/ul	98
26) Benzo(a)pyrene	23.181	252	18082	0.371	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.408	276	26965	0.369	ng/ul	97
28) Dibenzo(a,h)anthracene	25.425	278	21117	0.372	ng/ul	99
29) Benzo(g,h,i)perylene	26.059	276	21079	0.369	ng/ul	100

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

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