

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070124\
 Data File : BN032417.D
 Acq On : 03 Jul 2024 01:35
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4391

Quant Time: Jul 03 03:28:04 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN062724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 03 03:26:42 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.603	152	9404	0.400	ng/ul	0.00
4) Naphthalene-d8	10.382	136	28875	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.253	164	15340	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.008	188	27603	0.400	ng/ul	0.00
17) Chrysene-d12	21.209	240	23944	0.400	ng/ul	0.00
23) Perylene-d12	23.424	264	28993	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.148	96	3989	0.382	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.977	152	16683	0.385	ng/ul	0.00
18) Fluoranthene-d10	19.045	212	27770	0.385	ng/ul	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.181	88	4322	0.363	ng/ul	94
5) Naphthalene	10.431	128	31576	0.413	ng/ul	99
7) 2-Methylnaphthalene	12.054	142	19509	0.385	ng/ul	100
8) 1-Methylnaphthalene	12.274	142	20696	0.394	ng/ul	100
10) Acenaphthylene	13.971	152	28814	0.452	ng/ul	99
11) Acenaphthene	14.313	153	20692	0.418	ng/ul	98
12) Fluorene	15.313	166	21482	0.364	ng/ul	98
14) Pentachlorophenol	16.674	266	2819	0.451	ng/ul	100
15) Phenanthrene	17.050	178	30767	0.407	ng/ul	100
16) Anthracene	17.143	178	29156	0.432	ng/ul	99
19) Fluoranthene	19.073	202	35589	0.390	ng/ul	97
20) Pyrene	19.440	202	37103	0.389	ng/ul	99
21) Benzo(a)anthracene	21.191	228	36244	0.419	ng/ul	99
22) Chrysene	21.244	228	37280	0.393	ng/ul	98
24) Benzo(b)fluoranthene	22.757	252	40299	0.320	ng/ul	97
25) Benzo(k)fluoranthene	22.801	252	40021	0.327	ng/ul	98
26) Benzo(a)pyrene	23.327	252	39179	0.447	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.660	276	49054	0.357	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.666	278	39084	0.383	ng/ul	97
29) Benzo(g,h,i)perylene	26.344	276	40006	0.366	ng/ul	98

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

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