

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031824\
 Data File : BN030425.D
 Acq On : 18 Mar 2024 10:54
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4361

Quant Time: Mar 18 11:46:45 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN031524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 15 01:04:41 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.602	152	2078	0.400	ng/ul	0.00
4) Naphthalene-d8	10.380	136	4813	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.255	164	2706	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.014	188	5271	0.400	ng/ul	0.00
17) Chrysene-d12	21.220	240	4667	0.400	ng/ul	0.00
23) Perylene-d12	23.423	264	5371	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.155	96	1030	0.407	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.980	152	2594	0.390	ng/ul	0.00
18) Fluoranthene-d10	19.048	212	5646	0.393	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.189	88	1186	0.430	ng/ul	94
5) Naphthalene	10.429	128	5071	0.390	ng/ul	97
7) 2-Methylnaphthalene	12.057	142	2947	0.387	ng/ul#	100
8) 1-Methylnaphthalene	12.277	142	3144	0.390	ng/ul#	28
10) Acenaphthylene	13.973	152	5005	0.388	ng/ul	99
11) Acenaphthene	14.316	153	3696	0.391	ng/ul	100
12) Fluorene	15.315	166	4042	0.389	ng/ul	98
14) Pentachlorophenol	16.664	266	604	0.399	ng/ul#	100
15) Phenanthrene	17.056	178	5973	0.385	ng/ul	99
16) Anthracene	17.149	178	5332	0.376	ng/ul	99
19) Fluoranthene	19.080	202	7250	0.393	ng/ul	100
20) Pyrene	19.443	202	7805	0.396	ng/ul	98
21) Benzo(a)anthracene	21.205	228	5998	0.383	ng/ul	99
22) Chrysene	21.255	228	7089	0.394	ng/ul	98
24) Benzo(b)fluoranthene	22.765	252	6972	0.387	ng/ul	98
25) Benzo(k)fluoranthene	22.809	252	8542	0.390	ng/ul	99
26) Benzo(a)pyrene	23.333	252	7302	0.384	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.639	276	9343	0.372	ng/ul#	88
28) Dibenzo(a,h)anthracene	25.666	278	6965	0.363	ng/ul	98
29) Benzo(g,h,i)perylene	26.313	276	8570	0.382	ng/ul	99

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

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