

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN030224\
 Data File : BN030259.D
 Acq On : 04 Mar 2024 16:58
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 82 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4356

Quant Time: Mar 04 17:28:47 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN030224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 04 05:54:20 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.830	152	1412	0.400	ng/ul	0.00
4) Naphthalene-d8	10.628	136	3936	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.479	164	1899	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.235	188	3633	0.400	ng/ul #	0.00
17) Chrysene-d12	21.447	240	3191	0.400	ng/ul #	0.00
23) Perylene-d12	23.800	264	3848	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.307	96	1010	0.517	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.229	152	2527	0.497	ng/ul	0.00
18) Fluoranthene-d10	19.272	212	4797	0.461	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.340	88	1010	0.476	ng/ul	96
5) Naphthalene	10.678	128	5522	0.475	ng/ul	99
7) 2-Methylnaphthalene	12.300	142	3059	0.486	ng/ul	100
8) 1-Methylnaphthalene	12.515	142	3153	0.473	ng/ul	98
10) Acenaphthylene	14.197	152	4885	0.460	ng/ul	99
11) Acenaphthene	14.539	153	3551	0.456	ng/ul	98
12) Fluorene	15.538	166	3716	0.468	ng/ul	97
14) Pentachlorophenol	16.884	266	564	0.544	ng/ul	99
15) Phenanthrene	17.281	178	5495	0.451	ng/ul	100
16) Anthracene	17.374	178	5018	0.468	ng/ul	98
19) Fluoranthene	19.300	202	6295	0.444	ng/ul	98
20) Pyrene	19.667	202	6721	0.436	ng/ul#	94
21) Benzo(a)anthracene	21.430	228	5494	0.487	ng/ul	99
22) Chrysene	21.482	228	6020	0.444	ng/ul	99
24) Benzo(b)fluoranthene	23.084	252	6492	0.437	ng/ul	94
25) Benzo(k)fluoranthene	23.133	252	7444	0.425	ng/ul	94
26) Benzo(a)pyrene	23.698	252	6639	0.440	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	26.232	276	8736	0.453	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.259	278	6941	0.465	ng/ul	92
29) Benzo(g,h,i)perylene	26.970	276	7649	0.427	ng/ul#	93

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

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