

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031524\
 Data File : BN030411.D
 Acq On : 15 Mar 2024 09:44
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4358

Quant Time: Mar 15 11:06:02 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	7.604	152	2301	0.400	ng/ul	0.00
4) Naphthalene-d8	10.377	136	5742	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.254	164	3290	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.013	188	6629	0.400	ng/ul	0.00
17) Chrysene-d12	21.218	240	5930	0.400	ng/ul	0.00
23) Perylene-d12	23.428	264	6686	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.157	96	1114	0.397	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.983	152	3181	0.401	ng/ul	0.00
18) Fluoranthene-d10	19.046	212	7092	0.389	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.187	88	1189	0.389	ng/ul	94
5) Naphthalene	10.427	128	6022	0.388	ng/ul	98
7) 2-Methylnaphthalene	12.055	142	3582	0.395	ng/ul#	100
8) 1-Methylnaphthalene	12.275	142	3806	0.395	ng/ul#	28
10) Acenaphthylene	13.972	152	6083	0.388	ng/ul	99
11) Acenaphthene	14.319	153	4473	0.389	ng/ul	98
12) Fluorene	15.314	166	4994	0.395	ng/ul	98
14) Pentachlorophenol	16.662	266	715	0.375	ng/ul#	100
15) Phenanthrene	17.055	178	7529	0.385	ng/ul	97
16) Anthracene	17.148	178	6789	0.380	ng/ul	99
19) Fluoranthene	19.079	202	9034	0.385	ng/ul	99
20) Pyrene	19.446	202	9735	0.388	ng/ul	98
21) Benzo(a)anthracene	21.204	228	7746	0.389	ng/ul	100
22) Chrysene	21.253	228	8935	0.391	ng/ul	98
24) Benzo(b)fluoranthene	22.764	252	8685	0.387	ng/ul	96
25) Benzo(k)fluoranthene	22.808	252	10795	0.396	ng/ul	96
26) Benzo(a)pyrene	23.331	252	9111	0.385	ng/ul	92
27) Indeno(1,2,3-cd)pyrene	25.641	276	11491	0.367	ng/ul#	89
28) Dibenzo(a,h)anthracene	25.661	278	8835	0.370	ng/ul	95
29) Benzo(g,h,i)perylene	26.312	276	10406	0.372	ng/ul	98

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

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