

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012422\
 Data File : BM033835.D
 Acq On : 24 Jan 2022 13:18
 Operator : CG/JU
 Sample : SST0.846
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.8046

Quant Time: Jan 24 14:05:45 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM012422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 24 14:03:45 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	2023	0.400	ng/ul	0.00
4) Naphthalene-d8	10.756	136	6949	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.579	164	3547	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.316	188	7203	0.400	ng/ul	0.00
17) Chrysene-d12	21.484	240	6097	0.400	ng/ul #	0.00
23) Perylene-d12	23.885	264	6082	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.296	96	1684	0.781	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.343	152	7750	0.820	ng/ul	0.00
18) Fluoranthene-d10	19.337	212	14728	0.780	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.334	88	1762	0.750	ng/ul#	86
5) Naphthalene	10.805	128	15243	0.757	ng/ul#	88
7) 2-Methylnaphthalene	12.415	142	9977	0.772	ng/ul	100
8) 1-Methylnaphthalene	12.636	142	9750	0.768	ng/ul	100
10) Acenaphthylene	14.303	152	12379	0.755	ng/ul#	90
11) Acenaphthene	14.644	153	9944	0.760	ng/ul	94
12) Fluorene	15.626	166	11328	0.724	ng/ul#	97
14) Pentachlorophenol	16.972	266	2596	0.913	ng/ul	99
15) Phenanthrene	17.358	178	18007	0.772	ng/ul	95
16) Anthracene	17.451	178	16939	0.791	ng/ul#	91
19) Fluoranthene	19.368	202	21779	0.762	ng/ul	99
20) Pyrene	19.726	202	22103	0.764	ng/ul	95
21) Benzo(a)anthracene	21.467	228	19394	0.766	ng/ul	97
22) Chrysene	21.518	228	20464	0.790	ng/ul	97
24) Benzo(b)fluoranthene	23.150	252	21137	0.793	ng/ul	91
25) Benzo(k)fluoranthene	23.199	252	21651	0.828	ng/ul#	89
26) Benzo(a)pyrene	23.778	252	18852	0.814	ng/ul#	84
27) Indeno(1,2,3-cd)pyrene	26.392	276	24683	0.819	ng/ul#	85
28) Dibenzo(a,h)anthracene	26.412	278	19584	0.817	ng/ul#	88
29) Benzo(g,h,i)perylene	27.158	276	20992	0.810	ng/ul	94

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

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