

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040424\
 Data File : BM044899.D
 Acq On : 03 Apr 2024 20:02
 Operator : MA/JU
 Sample : SST0.818
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.8025

Quant Time: Apr 03 23:27:05 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM040424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 03 23:24:21 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.656	152	1911	0.400	ng/ul	0.00
4) Naphthalene-d8	10.435	136	5769	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.303	164	3042	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.065	188	6649	0.400	ng/ul	0.00
17) Chrysene-d12	21.271	240	6352	0.400	ng/ul	0.00
23) Perylene-d12	23.507	264	7809	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	1980	0.790	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.035	152	6348	0.766	ng/ul	-0.02
18) Fluoranthene-d10	19.104	212	13093	0.587	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.255	88	2022	0.790	ng/ul#	89
5) Naphthalene	10.484	128	12175	0.750	ng/ul	95
7) 2-Methylnaphthalene	12.112	142	7505	0.731	ng/ul	98
8) 1-Methylnaphthalene	12.327	142	8049	0.776	ng/ul	100
10) Acenaphthylene	14.021	152	12094	0.783	ng/ul	98
11) Acenaphthene	14.363	153	8715	0.781	ng/ul	99
12) Fluorene	15.367	166	9663	0.758	ng/ul	99
14) Pentachlorophenol	16.731	266	1184	0.738	ng/ul	97
15) Phenanthrene	17.107	178	14956	0.749	ng/ul	98
16) Anthracene	17.204	178	14468	0.772	ng/ul	100
19) Fluoranthene	19.132	202	18091	0.582	ng/ul	99
20) Pyrene	19.499	202	18920	0.594	ng/ul	100
21) Benzo(a)anthracene	21.256	228	17899	0.740	ng/ul	99
22) Chrysene	21.306	228	20456	0.753	ng/ul	100
24) Benzo(b)fluoranthene	22.832	252	21921	0.735	ng/ul	93
25) Benzo(k)fluoranthene	22.878	252	24138	0.748	ng/ul	95
26) Benzo(a)pyrene	23.410	252	21035	0.777	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	25.752	276	29310	0.941	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.775	278	23188	0.965	ng/ul	98
29) Benzo(g,h,i)perylene	26.429	276	24880	0.910	ng/ul	99

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

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