

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040424\
 Data File : BM044898.D
 Acq On : 03 Apr 2024 19:26
 Operator : MA/JU
 Sample : SST0.417
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.4024

Quant Time: Apr 03 23:26:53 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	1937	0.400	ng/ul	0.00
4) Naphthalene-d8	10.440	136	5585	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.307	164	3022	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.068	188	6981	0.400	ng/ul	0.00
17) Chrysene-d12	21.272	240	7240	0.400	ng/ul	0.00
23) Perylene-d12	23.505	264	9261	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.226	96	1025	0.404	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.052	152	3139	0.391	ng/ul	0.00
18) Fluoranthene-d10	19.103	212	7469	0.294	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.255	88	956	0.368	ng/ul	100
5) Naphthalene	10.495	128	6058	0.386	ng/ul	100
7) 2-Methylnaphthalene	12.129	142	3682	0.370	ng/ul	100
8) 1-Methylnaphthalene	12.338	142	3992	0.398	ng/ul	100
10) Acenaphthylene	14.029	152	6098	0.397	ng/ul	100
11) Acenaphthene	14.367	153	4466	0.403	ng/ul	100
12) Fluorene	15.375	166	5004	0.395	ng/ul	100
14) Pentachlorophenol	16.735	266	666	0.396	ng/ul	96
15) Phenanthrene	17.111	178	7978	0.381	ng/ul	100
16) Anthracene	17.212	178	7894	0.401	ng/ul	100
19) Fluoranthene	19.136	202	10308	0.291	ng/ul	100
20) Pyrene	19.503	202	10959	0.302	ng/ul	100
21) Benzo(a)anthracene	21.260	228	10280	0.373	ng/ul	100
22) Chrysene	21.310	228	12606	0.407	ng/ul	100
24) Benzo(b)fluoranthene	22.835	252	13422	0.380	ng/ul	100
25) Benzo(k)fluoranthene	22.882	252	15069	0.394	ng/ul	100
26) Benzo(a)pyrene	23.414	252	12891	0.402	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.756	276	17718	0.480	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.783	278	14062	0.494	ng/ul	100
29) Benzo(g,h,i)perylene	26.440	276	15133	0.467	ng/ul	100

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

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