

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070524\
 Data File : BM046379.D
 Acq On : 05 Jul 2024 12:49
 Operator : MA/JU
 Sample : SST0.264
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.2030

Quant Time: Jul 05 14:56:20 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM070524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 05 14:54:52 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.538	152	4899	0.400	ng/ul	0.00
4) Naphthalene-d8	10.303	136	13322	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.182	164	7665	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.933	188	15326	0.400	ng/ul	0.00
17) Chrysene-d12	21.146	240	9889	0.400	ng/ul	0.00
23) Perylene-d12	23.306	264	10281	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.133	96	1047	0.165	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.903	152	3942	0.220	ng/ul	0.00
18) Fluoranthene-d10	18.973	212	7076	0.241	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.163	88	1234	0.177	ng/ul#	88
5) Naphthalene	10.352	128	7047	0.195	ng/ul	99
7) 2-Methylnaphthalene	11.980	142	4776	0.227	ng/ul	99
8) 1-Methylnaphthalene	12.200	142	4899	0.215	ng/ul	100
10) Acenaphthylene	13.899	152	7750	0.216	ng/ul	99
11) Acenaphthene	14.246	153	5036	0.198	ng/ul	100
12) Fluorene	15.241	166	5930	0.213	ng/ul	98
14) Pentachlorophenol	16.583	266	777	0.175	ng/ul	96
15) Phenanthrene	16.975	178	9092	0.213	ng/ul	99
16) Anthracene	17.068	178	8776	0.261	ng/ul	98
19) Fluoranthene	19.006	202	10144	0.239	ng/ul	99
20) Pyrene	19.368	202	10436	0.230	ng/ul	99
21) Benzo(a)anthracene	21.132	228	9002	0.236	ng/ul	100
22) Chrysene	21.181	228	8946	0.192	ng/ul	100
24) Benzo(b)fluoranthene	22.666	252	8979	0.242	ng/ul	98
25) Benzo(k)fluoranthene	22.710	252	8776	0.204	ng/ul	97
26) Benzo(a)pyrene	23.212	252	6848	0.205	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	25.451	276	9359	0.175	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.477	278	7209	0.180	ng/ul	97
29) Benzo(g,h,i)perylene	26.095	276	7466	0.164	ng/ul	97

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

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