

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090321\
 Data File : BM032037.D
 Acq On : 03 Sep 2021 12:32
 Operator : CG/JU
 Sample : SST0.122
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.1022

Quant Time: Sep 03 14:41:35 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM090321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 03 14:29:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.452	152	1341	0.400	ng/ul	0.00
4) Naphthalene-d8	10.212	136	4588	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.098	164	2855	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.864	188	5836	0.400	ng/ul	0.00
17) Chrysene-d12	21.072	240	4976	0.400	ng/ul	0.00
23) Perylene-d12	23.193	264	4839	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.071	96	158	0.106	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.813	152	752	0.097	ng/ul	0.00
18) Fluoranthene-d10	18.902	212	1554	0.094	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.102	88	233	0.155	ng/ul#	42
5) Naphthalene	10.262	128	1301	0.095	ng/ul#	89
7) 2-Methylnaphthalene	11.885	142	879	0.094	ng/ul	98
8) 1-Methylnaphthalene	12.110	142	886	0.096	ng/ul	99
10) Acenaphthylene	13.824	152	1237	0.102	ng/ul#	89
11) Acenaphthene	14.163	153	1024	0.100	ng/ul	96
12) Fluorene	15.164	166	1220	0.104	ng/ul#	96
14) Pentachlorophenol	16.530	266	196	0.105	ng/ul	93
15) Phenanthrene	16.905	178	1977	0.107	ng/ul	95
16) Anthracene	16.999	178	1602	0.092	ng/ul	94
19) Fluoranthene	18.932	202	2413	0.114	ng/ul#	87
20) Pyrene	19.294	202	2649	0.123	ng/ul#	88
21) Benzo(a)anthracene	21.057	228	2095	0.104	ng/ul	98
22) Chrysene	21.108	228	2384	0.115	ng/ul	96
24) Benzo(b)fluoranthene	22.609	252	2533	0.116	ng/ul	69
25) Benzo(k)fluoranthene	22.609	252	2533	0.112	ng/ul#	70
26) Benzo(a)pyrene	23.103	252	2111	0.111	ng/ul#	61
27) Indeno(1,2,3-cd)pyrene	25.276	276	2677	0.108	ng/ul#	93
28) Dibenzo(a,h)anthracene	25.291	278	2188	0.110	ng/ul#	76
29) Benzo(g,h,i)perylene	25.911	276	2478	0.118	ng/ul#	89

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

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