

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090321\
 Data File : BM032040.D
 Acq On : 03 Sep 2021 14:33
 Operator : CG/JU
 Sample : SST0.825
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.8025

Quant Time: Sep 03 15:49:40 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	7.459	152	1158	0.400	ng/ul	0.00
4) Naphthalene-d8	10.212	136	3946	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.102	164	2494	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.862	188	4812	0.400	ng/ul	0.00
17) Chrysene-d12	21.069	240	3673	0.400	ng/ul	0.00
23) Perylene-d12	23.192	264	3445	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.068	96	1037	0.806	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.813	152	4903	0.737	ng/ul	0.00
18) Fluoranthene-d10	18.900	212	8433	0.688	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.099	88	994	0.765	ng/ul#	89
5) Naphthalene	10.266	128	8217	0.699	ng/ul	96
7) 2-Methylnaphthalene	11.894	142	5841	0.727	ng/ul	100
8) 1-Methylnaphthalene	12.110	142	5736	0.723	ng/ul	99
10) Acenaphthylene	13.824	152	7882	0.742	ng/ul	99
11) Acenaphthene	14.167	153	6302	0.705	ng/ul	99
12) Fluorene	15.164	166	7300	0.710	ng/ul	99
14) Pentachlorophenol	16.532	266	979	0.635	ng/ul	93
15) Phenanthrene	16.904	178	11106	0.727	ng/ul	97
16) Anthracene	17.001	178	10065	0.704	ng/ul	97
19) Fluoranthene	18.930	202	12132	0.774	ng/ul	98
20) Pyrene	19.296	202	12109	0.761	ng/ul	97
21) Benzo(a)anthracene	21.057	228	9973	0.672	ng/ul	97
22) Chrysene	21.107	228	11340	0.739	ng/ul	99
24) Benzo(b)fluoranthene	22.562	252	12091	0.779	ng/ul	91
25) Benzo(k)fluoranthene	22.604	252	12021	0.750	ng/ul#	89
26) Benzo(a)pyrene	23.100	252	10475	0.771	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	25.264	276	14844	0.845	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.276	278	12041	0.850	ng/ul	90
29) Benzo(g,h,i)perylene	25.896	276	13186	0.880	ng/ul	98

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

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