

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071321\
 Data File : BM030926.D
 Acq On : 13 Jul 2021 12:14
 Operator : CG/JU
 Sample : SST0.804
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.8004

Quant Time: Jul 13 12:47:04 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM071321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 13 12:12:03 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.673	152	775	0.400	ng/ul	0.00
4) Naphthalene-d8	10.442	136	2546	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.303	164	1664	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.048	188	3487	0.400	ng/ul	0.00
17) Chrysene-d12	21.236	240	3051	0.400	ng/ul	0.00
23) Perylene-d12	23.433	264	2941	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.241	96	641	0.759	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.034	152	3636	0.909	ng/ul	0.00
18) Fluoranthene-d10	19.077	212	8379	0.958	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.272	88	725	0.831	ng/ul	91
5) Naphthalene	10.491	128	6090	0.839	ng/ul	96
7) 2-Methylnaphthalene	12.106	142	4410	0.903	ng/ul	99
8) 1-Methylnaphthalene	12.331	142	4271	0.899	ng/ul	99
10) Acenaphthylene	14.019	152	6466	0.870	ng/ul	97
11) Acenaphthene	14.364	153	4685	0.778	ng/ul	98
12) Fluorene	15.353	166	5603	0.844	ng/ul	100
14) Pentachlorophenol	16.697	266	1008	0.886	ng/ul	99
15) Phenanthrene	17.089	178	8938	0.762	ng/ul	99
16) Anthracene	17.179	178	8549	0.833	ng/ul	98
19) Fluoranthene	19.108	202	10771	0.800	ng/ul	98
20) Pyrene	19.470	202	10896	0.805	ng/ul	100
21) Benzo(a)anthracene	21.222	228	10648	0.876	ng/ul	99
22) Chrysene	21.272	228	10965	0.805	ng/ul	98
24) Benzo(b)fluoranthene	22.776	252	11878	0.912	ng/ul	92
25) Benzo(k)fluoranthene	22.822	252	11589	0.863	ng/ul#	92
26) Benzo(a)pyrene	23.336	252	10157	0.917	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	25.618	276	14030	0.972	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.638	278	11387	0.983	ng/ul	93
29) Benzo(g,h,i)perylene	26.282	276	11911	0.950	ng/ul	98

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

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