

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071521\
 Data File : BM030980.D
 Acq On : 15 Jul 2021 14:26
 Operator : CG/JU
 Sample : PB137577BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS577

Quant Time: Jul 15 15:01:36 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 : Thu Jul 15 13:02:53 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.652	152	985	0.400	ng/ul	0.00
4) Naphthalene-d8	10.419	136	3181	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.284	164	1898	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.030	188	4128	0.400	ng/ul	0.00
17) Chrysene-d12	21.222	240	3638	0.400	ng/ul	0.00
23) Perylene-d12	23.409	264	3563	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.227	96	757	0.747	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.016	152	1776	0.321	ng/ul	0.00
18) Fluoranthene-d10	19.062	212	4324	0.353	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.258	88	1961	1.735	ng/ul#	82
5) Naphthalene	10.473	128	3016	0.318	ng/ul	99
7) 2-Methylnaphthalene	12.088	142	2077	0.310	ng/ul	99
8) 1-Methylnaphthalene	12.313	142	1908	0.291	ng/ul	100
10) Acenaphthylene	14.004	152	2470	0.266	ng/ul	97
11) Acenaphthene	14.349	153	2150	0.320	ng/ul	95
12) Fluorene	15.338	166	2529	0.317	ng/ul	99
14) Pentachlorophenol	16.680	266	1354	0.929	ng/ul	99
15) Phenanthrene	17.072	178	4261	0.329	ng/ul	99
16) Anthracene	17.162	178	3928	0.316	ng/ul	99
19) Fluoranthene	19.092	202	5282	0.330	ng/ul	99
20) Pyrene	19.454	202	5461	0.337	ng/ul	98
21) Benzo(a)anthracene	21.207	228	5285	0.336	ng/ul	100
22) Chrysene	21.258	228	5556	0.342	ng/ul	99
24) Benzo(b)fluoranthene	22.754	252	5909	0.343	ng/ul	98
25) Benzo(k)fluoranthene	22.800	252	5937	0.343	ng/ul	97
26) Benzo(a)pyrene	23.312	252	4972	0.332	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.582	276	7164	0.353	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.601	278	5732	0.349	ng/ul	96
29) Benzo(g,h,i)perylene	26.243	276	6079	0.353	ng/ul	99

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

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