

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071321\
 Data File : BM030965.D
 Acq On : 14 Jul 2021 18:09
 Operator : CG/JU
 Sample : PB137577BS
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS577

Quant Time: Jul 14 18:39:27 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 14:01:01 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	1043	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.428	136	3469	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.292	164	2151	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.041	188	4354	0.400	ng/ul	-0.01
17) Chrysene-d12	21.227	240	3390	0.400	ng/ul	0.00
23) Perylene-d12	23.416	264	3220	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.237	96	750	0.699	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.025	152	1950	0.323	ng/ul	0.00
18) Fluoranthene-d10	19.069	212	4333	0.380	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.269	88	2008	1.678	ng/ul#	80
5) Naphthalene	10.482	128	3283	0.317	ng/ul	98
7) 2-Methylnaphthalene	12.097	142	2274	0.311	ng/ul	100
8) 1-Methylnaphthalene	12.317	142	2092	0.293	ng/ul	99
10) Acenaphthylene	14.011	152	2924	0.277	ng/ul	99
11) Acenaphthene	14.356	153	2378	0.313	ng/ul	98
12) Fluorene	15.346	166	2770	0.306	ng/ul	99
14) Pentachlorophenol	16.687	266	1380	0.897	ng/ul	100
15) Phenanthrene	17.079	178	4469	0.327	ng/ul	99
16) Anthracene	17.172	178	4114	0.314	ng/ul	100
19) Fluoranthene	19.100	202	5276	0.353	ng/ul	99
20) Pyrene	19.462	202	5347	0.354	ng/ul	99
21) Benzo(a)anthracene	21.210	228	4892	0.334	ng/ul	99
22) Chrysene	21.260	228	5213	0.344	ng/ul	99
24) Benzo(b)fluoranthene	22.759	252	5466	0.351	ng/ul	98
25) Benzo(k)fluoranthene	22.803	252	5553	0.355	ng/ul	99
26) Benzo(a)pyrene	23.319	252	4495	0.332	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.594	276	6539	0.356	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.608	278	5198	0.350	ng/ul	98
29) Benzo(g,h,i)perylene	26.255	276	5465	0.351	ng/ul	99

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

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