

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012422\
 Data File : BM033846.D
 Acq On : 24 Jan 2022 20:12
 Operator : CG/JU
 Sample : PB142204BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS204

Quant Time: Jan 25 02:21:01 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM012422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 24 15:20:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.943	152	2205	0.400	ng/ul	0.00
4) Naphthalene-d8	10.756	136	7837	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.580	164	4018	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.320	188	8298	0.400	ng/ul	0.00
17) Chrysene-d12	21.480	240	6955	0.400	ng/ul #	0.00
23) Perylene-d12	23.880	264	6909	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.296	96	1459	0.633	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.344	152	3817	0.353	ng/ul	0.00
18) Fluoranthene-d10	19.338	212	8207	0.383	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.334	88	3745	1.552	ng/ul#	80
5) Naphthalene	10.805	128	7194	0.334	ng/ul#	89
7) 2-Methylnaphthalene	12.416	142	4552	0.325	ng/ul	99
8) 1-Methylnaphthalene	12.636	142	4442	0.324	ng/ul	100
10) Acenaphthylene	14.303	152	6668	0.371	ng/ul#	92
11) Acenaphthene	14.640	153	4830	0.336	ng/ul	97
12) Fluorene	15.626	166	5406	0.330	ng/ul	97
14) Pentachlorophenol	16.972	266	2422	0.635	ng/ul	97
15) Phenanthrene	17.358	178	9066	0.343	ng/ul	95
16) Anthracene	17.451	178	8576	0.347	ng/ul	93
19) Fluoranthene	19.365	202	11697	0.369	ng/ul	98
20) Pyrene	19.727	202	12140	0.374	ng/ul	95
21) Benzo(a)anthracene	21.465	228	10354	0.364	ng/ul	98
22) Chrysene	21.516	228	11132	0.369	ng/ul	99
24) Benzo(b)fluoranthene	23.148	252	11085	0.357	ng/ul	93
25) Benzo(k)fluoranthene	23.195	252	11226	0.361	ng/ul#	92
26) Benzo(a)pyrene	23.774	252	10794	0.401	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	26.390	276	12410	0.347	ng/ul#	87
28) Dibenzo(a,h)anthracene	26.410	278	9939	0.352	ng/ul	95
29) Benzo(g,h,i)perylene	27.153	276	10622	0.350	ng/ul#	93

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

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