

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111921\
 Data File : BM033185.D
 Acq On : 19 Nov 2021 20:41
 Operator : CG/JU
 Sample : PB140811BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS811

Quant Time: Nov 20 01:05:19 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM111921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 19 15:41:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	2438	0.400	ng/ul	0.00
4) Naphthalene-d8	10.760	136	6866	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.579	164	4316	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.316	188	9802	0.400	ng/ul	0.00
17) Chrysene-d12	21.472	240	7946	0.400	ng/ul	0.00
23) Perylene-d12	23.812	264	6132	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.420	96	1343	0.586	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.339	152	3097	0.326	ng/ul	0.00
18) Fluoranthene-d10	19.330	212	7893	0.343	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.451	88	4058	1.574	ng/ul	93
5) Naphthalene	10.810	128	6582	0.309	ng/ul	99
7) 2-Methylnaphthalene	12.415	142	4478	0.325	ng/ul	96
8) 1-Methylnaphthalene	12.631	142	4277	0.314	ng/ul	99
10) Acenaphthylene	14.303	152	6339	0.314	ng/ul#	99
11) Acenaphthene	14.640	153	5245	0.317	ng/ul	98
12) Fluorene	15.626	166	6099	0.327	ng/ul	99
14) Pentachlorophenol	16.962	266	1648	0.504	ng/ul	98
15) Phenanthrene	17.357	178	10015	0.282	ng/ul	99
16) Anthracene	17.451	178	8771	0.305	ng/ul	100
19) Fluoranthene	19.360	202	11657	0.305	ng/ul	98
20) Pyrene	19.722	202	11735	0.309	ng/ul	96
21) Benzo(a)anthracene	21.453	228	9550	0.337	ng/ul	99
22) Chrysene	21.508	228	10528	0.337	ng/ul	98
24) Benzo(b)fluoranthene	23.102	252	9923	0.357	ng/ul	99
25) Benzo(k)fluoranthene	23.148	252	9559	0.349	ng/ul	98
26) Benzo(a)pyrene	23.708	252	7790	0.339	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.208	276	9429	0.325	ng/ul#	89
28) Dibenzo(a,h)anthracene	26.222	278	7487	0.323	ng/ul	96
29) Benzo(g,h,i)perylene	26.942	276	8424	0.323	ng/ul	95

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

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