

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111921\
 Data File : BM033186.D
 Acq On : 19 Nov 2021 21:17
 Operator : CG/JU
 Sample : PB140818BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS818

Quant Time: Nov 20 01:05:32 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	1817	0.400	ng/ul	0.00
4) Naphthalene-d8	10.760	136	5138	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.580	164	3252	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.316	188	7489	0.400	ng/ul	0.00
17) Chrysene-d12	21.470	240	6903	0.400	ng/ul	0.00
23) Perylene-d12	23.813	264	5824	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.416	96	1227	0.718	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.339	152	2729	0.383	ng/ul	0.00
18) Fluoranthene-d10	19.331	212	7400	0.370	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.451	88	3494	1.818	ng/ul	93
5) Naphthalene	10.810	128	5675	0.356	ng/ul	99
7) 2-Methylnaphthalene	12.415	142	3843	0.373	ng/ul	96
8) 1-Methylnaphthalene	12.631	142	3733	0.366	ng/ul	99
10) Acenaphthylene	14.303	152	5458	0.359	ng/ul#	97
11) Acenaphthene	14.644	153	4533	0.363	ng/ul	99
12) Fluorene	15.623	166	5326	0.378	ng/ul	98
14) Pentachlorophenol	16.962	266	1390	0.556	ng/ul	98
15) Phenanthrene	17.358	178	8817	0.325	ng/ul	99
16) Anthracene	17.448	178	7731	0.352	ng/ul	98
19) Fluoranthene	19.361	202	10510	0.316	ng/ul	97
20) Pyrene	19.723	202	10738	0.326	ng/ul	96
21) Benzo(a)anthracene	21.453	228	9124	0.370	ng/ul	99
22) Chrysene	21.507	228	10059	0.371	ng/ul	99
24) Benzo(b)fluoranthene	23.098	252	10268	0.389	ng/ul	99
25) Benzo(k)fluoranthene	23.146	252	9907	0.381	ng/ul	99
26) Benzo(a)pyrene	23.706	252	8119	0.371	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.209	276	10021	0.364	ng/ul#	87
28) Dibenzo(a,h)anthracene	26.221	278	7907	0.359	ng/ul	96
29) Benzo(g,h,i)perylene	26.943	276	8905	0.360	ng/ul	95

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

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