

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111321\
 Data File : BM033073.D
 Acq On : 14 Nov 2021 10:59
 Operator : CG/JU
 Sample : PB140482BS
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS482

Quant Time: Nov 15 03:45:14 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM110921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 15 03:11:31 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.438	152	1425	0.400	ng/ul	0.00
4) Naphthalene-d8	10.203	136	5940	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.091	164	3747	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.834	188	7428	0.400	ng/ul	0.00
17) Chrysene-d12	21.030	240	5127	0.400	ng/ul	0.00
23) Perylene-d12	23.134	264	4114	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.843	96	1364	0.750	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	11.804	152	3088	0.366	ng/ul	0.00
18) Fluoranthene-d10	18.869	212	6470	0.416	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.878	88	3628	1.941	ng/ul#	90
5) Naphthalene	10.253	128	6212	0.360	ng/ul	100
7) 2-Methylnaphthalene	11.881	142	4145	0.347	ng/ul	99
8) 1-Methylnaphthalene	12.101	142	3990	0.341	ng/ul	100
10) Acenaphthylene	13.810	152	5448	0.334	ng/ul	99
11) Acenaphthene	14.152	153	4879	0.343	ng/ul	99
12) Fluorene	15.145	166	5407	0.322	ng/ul	99
14) Pentachlorophenol	16.511	266	963	0.498	ng/ul	97
15) Phenanthrene	16.876	178	8233	0.334	ng/ul	99
16) Anthracene	16.969	178	6886	0.317	ng/ul	99
19) Fluoranthene	18.900	202	9132	0.380	ng/ul	100
20) Pyrene	19.262	202	9283	0.371	ng/ul	100
21) Benzo(a)anthracene	21.016	228	6397	0.338	ng/ul	99
22) Chrysene	21.067	228	7542	0.353	ng/ul	100
24) Benzo(b)fluoranthene	22.514	252	6587	0.349	ng/ul	94
25) Benzo(k)fluoranthene	22.555	252	7260	0.359	ng/ul	92
26) Benzo(a)pyrene	23.045	252	5544	0.356	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	25.187	276	6924	0.387	ng/ul#	91
28) Dibenzo(a,h)anthracene	25.196	278	5445	0.384	ng/ul	95
29) Benzo(g,h,i)perylene	25.817	276	6325	0.407	ng/ul	99

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

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