

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072921\
 Data File : BM031331.D
 Acq On : 01 Aug 2021 03:26
 Operator : CG/JU
 Sample : PB137930BS
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS930

Quant Time: Aug 02 01:15:07 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 30 14:16:52 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.604	152	1216	0.400	ng/ul	0.00
4) Naphthalene-d8	10.369	136	4628	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.235	164	2680	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.985	188	5021	0.400	ng/ul	0.00
17) Chrysene-d12	21.173	240	3634	0.400	ng/ul	0.00
23) Perylene-d12	23.336	264	3446	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.196	96	890	0.659	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.961	152	2860	0.367	ng/ul	0.00
18) Fluoranthene-d10	19.016	212	5021	0.414	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.227	88	2328	1.706	ng/ul#	80
5) Naphthalene	10.418	128	4948	0.359	ng/ul	99
7) 2-Methylnaphthalene	12.037	142	3399	0.361	ng/ul	100
8) 1-Methylnaphthalene	12.258	142	3170	0.341	ng/ul	100
10) Acenaphthylene	13.958	152	4316	0.378	ng/ul	100
11) Acenaphthene	14.299	153	3443	0.358	ng/ul	99
12) Fluorene	15.293	166	3809	0.345	ng/ul	99
14) Pentachlorophenol	16.638	266	955	0.594	ng/ul	99
15) Phenanthrene	17.026	178	5830	0.366	ng/ul	98
16) Anthracene	17.117	178	5434	0.364	ng/ul	100
19) Fluoranthene	19.046	202	6248	0.403	ng/ul	97
20) Pyrene	19.408	202	6242	0.397	ng/ul	96
21) Benzo(a)anthracene	21.156	228	5669	0.386	ng/ul	99
22) Chrysene	21.207	228	5892	0.388	ng/ul	99
24) Benzo(b)fluoranthene	22.691	252	5933	0.382	ng/ul	96
25) Benzo(k)fluoranthene	22.732	252	6023	0.375	ng/ul	97
26) Benzo(a)pyrene	23.241	252	5227	0.384	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	25.477	276	6653	0.379	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.492	278	5429	0.383	ng/ul	98
29) Benzo(g,h,i)perylene	26.131	276	5713	0.381	ng/ul	97

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

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