

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081221\
 Data File : BM031670.D
 Acq On : 12 Aug 2021 10:39
 Operator : CG/JU
 Sample : PB138138BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS138

Quant Time: Aug 12 11:28:17 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 : Thu Aug 12 11:27:44 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.573	152	930	0.400	ng/ul	0.00
4) Naphthalene-d8	10.334	136	3306	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.209	164	1678	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.964	188	3140	0.400	ng/ul #	0.00
17) Chrysene-d12	21.156	240	2458	0.400	ng/ul #	0.00
23) Perylene-d12	23.317	264	2386	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.186	96	745	0.721	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.935	152	1946	0.349	ng/ul	0.00
18) Fluoranthene-d10	18.997	212	3235	0.394	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.217	88	1858	1.781	ng/ul#	78
5) Naphthalene	10.383	128	3929	0.399	ng/ul	98
7) 2-Methylnaphthalene	12.011	142	2526	0.375	ng/ul	100
8) 1-Methylnaphthalene	12.227	142	2360	0.355	ng/ul	98
10) Acenaphthylene	13.932	152	3019	0.422	ng/ul	99
11) Acenaphthene	14.273	153	2419	0.402	ng/ul	99
12) Fluorene	15.270	166	2607	0.377	ng/ul	99
14) Pentachlorophenol	16.628	266	707	0.703	ng/ul	99
15) Phenanthrene	17.006	178	3891	0.390	ng/ul	99
16) Anthracene	17.096	178	3751	0.402	ng/ul	97
19) Fluoranthene	19.028	202	4435	0.423	ng/ul#	94
20) Pyrene	19.390	202	4602	0.432	ng/ul#	95
21) Benzo(a)anthracene	21.144	228	4048	0.408	ng/ul	96
22) Chrysene	21.195	228	4294	0.418	ng/ul	97
24) Benzo(b)fluoranthene	22.672	252	4253	0.396	ng/ul	89
25) Benzo(k)fluoranthene	22.718	252	4402	0.396	ng/ul#	90
26) Benzo(a)pyrene	23.225	252	3775	0.401	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	25.456	276	4851	0.399	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.468	278	3756	0.383	ng/ul	91
29) Benzo(g,h,i)perylene	26.105	276	4226	0.407	ng/ul#	92

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

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