

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012623\
 Data File : BM038546.D
 Acq On : 26 Jan 2023 13:11
 Operator : CG/JU
 Sample : PB150439BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS439

Quant Time: Jan 27 00:43:30 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM012623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 27 00:18:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.971	152	1241	0.400	ng/ul	0.00
4) Naphthalene-d8	10.779	136	5103	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.601	164	2952	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.345	188	6078	0.400	ng/ul	0.00
17) Chrysene-d12	21.518	240	5369	0.400	ng/ul	0.00
23) Perylene-d12	23.933	264	5176	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.343	96	1215	0.746	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.363	152	2186	0.276	ng/ul	0.00
18) Fluoranthene-d10	19.366	212	4987	0.325	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.376	88	3136	1.896	ng/ul#	23
5) Naphthalene	10.829	128	3839	0.231	ng/ul	99
7) 2-Methylnaphthalene	12.434	142	2395	0.273	ng/ul	97
8) 1-Methylnaphthalene	12.654	142	2493	0.276	ng/ul	98
10) Acenaphthylene	14.324	152	3632	0.279	ng/ul	99
11) Acenaphthene	14.666	153	2793	0.282	ng/ul	100
12) Fluorene	15.651	166	3212	0.259	ng/ul	100
14) Pentachlorophenol	16.999	266	1149	0.464	ng/ul	99
15) Phenanthrene	17.388	178	5087	0.259	ng/ul	98
16) Anthracene	17.481	178	4649	0.278	ng/ul	97
19) Fluoranthene	19.399	202	6222	0.325	ng/ul	98
20) Pyrene	19.761	202	6544	0.318	ng/ul	99
21) Benzo(a)anthracene	21.501	228	5662	0.297	ng/ul	100
22) Chrysene	21.556	228	5839	0.305	ng/ul	99
24) Benzo(b)fluoranthene	23.193	252	6507	0.334	ng/ul	94
25) Benzo(k)fluoranthene	23.243	252	6391	0.338	ng/ul#	96
26) Benzo(a)pyrene	23.825	252	5768	0.330	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	26.441	276	7403	0.301	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.465	278	5895	0.301	ng/ul	95
29) Benzo(g,h,i)perylene	27.223	276	6655	0.311	ng/ul	99

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

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