

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012623\
 Data File : BM038601.D
 Acq On : 28 Jan 2023 00:52
 Operator : CG/JU
 Sample : PB150411BS
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS411

Quant Time: Jan 28 01:34:34 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	7.963	152	1103	0.400	ng/ul	0.00
4) Naphthalene-d8	10.768	136	3119	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.592	164	1933	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.337	188	4099	0.400	ng/ul	0.00
17) Chrysene-d12	21.510	240	5118	0.400	ng/ul	0.00
23) Perylene-d12	23.918	264	6807	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.343	96	701	0.484	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.357	152	1382	0.286	ng/ul	0.00
18) Fluoranthene-d10	19.362	212	3512	0.240	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.376	88	1716	1.167	ng/ul#	18
5) Naphthalene	10.818	128	2265	0.223	ng/ul	97
7) 2-Methylnaphthalene	12.429	142	1513	0.282	ng/ul	96
8) 1-Methylnaphthalene	12.649	142	1563	0.283	ng/ul	100
10) Acenaphthylene	14.319	152	2427	0.285	ng/ul	96
11) Acenaphthene	14.657	153	1775	0.273	ng/ul	98
12) Fluorene	15.642	166	2075	0.256	ng/ul	99
14) Pentachlorophenol	16.995	266	939	0.562	ng/ul	98
15) Phenanthrene	17.379	178	3355	0.253	ng/ul	96
16) Anthracene	17.472	178	3344	0.296	ng/ul	96
19) Fluoranthene	19.390	202	4507	0.247	ng/ul	98
20) Pyrene	19.752	202	4904	0.250	ng/ul	99
21) Benzo(a)anthracene	21.495	228	5237	0.288	ng/ul	97
22) Chrysene	21.548	228	5318	0.291	ng/ul	98
24) Benzo(b)fluoranthene	23.179	252	6886	0.269	ng/ul	95
25) Benzo(k)fluoranthene	23.228	252	6905	0.278	ng/ul	96
26) Benzo(a)pyrene	23.813	252	6340	0.276	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.424	276	9615	0.297	ng/ul#	65
28) Dibenzo(a,h)anthracene	26.441	278	7770	0.302	ng/ul	96
29) Benzo(g,h,i)perylene	27.196	276	8354	0.297	ng/ul	98

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

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