

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012623\
 Data File : BM038575.D
 Acq On : 27 Jan 2023 08:16
 Operator : CG/JU
 Sample : PB150409BS
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS409

Quant Time: Jan 27 21:58:33 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	7.967	152	1042	0.400	ng/ul	0.00
4) Naphthalene-d8	10.774	136	2534	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.597	164	1606	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.341	188	3057	0.400	ng/ul	0.00
17) Chrysene-d12	21.515	240	2810	0.400	ng/ul	0.00
23) Perylene-d12	23.927	264	4093	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.343	96	1009	0.737	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.357	152	1256	0.319	ng/ul	0.00
18) Fluoranthene-d10	19.362	212	2747	0.342	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.376	88	2565	1.847	ng/ul#	22
5) Naphthalene	10.823	128	2204	0.267	ng/ul	96
7) 2-Methylnaphthalene	12.434	142	1377	0.316	ng/ul	97
8) 1-Methylnaphthalene	12.649	142	1444	0.321	ng/ul	100
10) Acenaphthylene	14.319	152	2364	0.334	ng/ul	96
11) Acenaphthene	14.657	153	1751	0.324	ng/ul	98
12) Fluorene	15.642	166	1956	0.290	ng/ul	98
14) Pentachlorophenol	16.995	266	753	0.605	ng/ul	100
15) Phenanthrene	17.379	178	2921	0.296	ng/ul	96
16) Anthracene	17.476	178	2841	0.337	ng/ul	97
19) Fluoranthene	19.394	202	3540	0.353	ng/ul	99
20) Pyrene	19.757	202	3706	0.344	ng/ul	98
21) Benzo(a)anthracene	21.498	228	3491	0.350	ng/ul	99
22) Chrysene	21.553	228	3576	0.357	ng/ul	98
24) Benzo(b)fluoranthene	23.187	252	4916	0.319	ng/ul	96
25) Benzo(k)fluoranthene	23.237	252	4957	0.331	ng/ul	96
26) Benzo(a)pyrene	23.816	252	4762	0.344	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.434	276	8025	0.412	ng/ul	94
28) Dibenzo(a,h)anthracene	26.448	278	6478	0.418	ng/ul	98
29) Benzo(g,h,i)perylene	27.209	276	7248	0.429	ng/ul	97

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

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