

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012623\
 Data File : BM038557.D
 Acq On : 26 Jan 2023 20:40
 Operator : CG/JU
 Sample : PB150403BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS403

Quant Time: Jan 27 00:45:00 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	1001	0.400	ng/u1	0.00
4) Naphthalene-d8	10.779	136	2405	0.400	ng/u1 #	0.00
9) Acenaphthene-d10	14.601	164	1336	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.345	188	2772	0.400	ng/u1	0.00
17) Chrysene-d12	21.515	240	3099	0.400	ng/u1	0.00
23) Perylene-d12	23.927	264	4492	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.343	96	879	0.669	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.363	152	1157	0.310	ng/u1	0.00
18) Fluoranthene-d10	19.366	212	2702	0.305	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.381	88	2110	1.582	ng/u1#	22
5) Naphthalene	10.828	128	1929	0.247	ng/u1	94
7) 2-Methylnaphthalene	12.434	142	1142	0.277	ng/u1	99
8) 1-Methylnaphthalene	12.654	142	1262	0.296	ng/u1	99
10) Acenaphthylene	14.324	152	1787	0.303	ng/u1	96
11) Acenaphthene	14.661	153	1339	0.298	ng/u1	99
12) Fluorene	15.647	166	1513	0.270	ng/u1	99
14) Pentachlorophenol	16.999	266	632	0.560	ng/u1	99
15) Phenanthrene	17.383	178	2375	0.265	ng/u1	93
16) Anthracene	17.481	178	2288	0.299	ng/u1	94
19) Fluoranthene	19.394	202	3090	0.280	ng/u1	98
20) Pyrene	19.757	202	3281	0.276	ng/u1	99
21) Benzo(a)anthracene	21.501	228	3303	0.300	ng/u1	99
22) Chrysene	21.553	228	3420	0.309	ng/u1	96
24) Benzo(b)fluoranthene	23.190	252	4694	0.278	ng/u1	96
25) Benzo(k)fluoranthene	23.240	252	4843	0.295	ng/u1	97
26) Benzo(a)pyrene	23.822	252	4677	0.308	ng/u1	94
27) Indeno(1,2,3-cd)pyrene	26.441	276	7539	0.353	ng/u1#	95
28) Dibenzo(a,h)anthracene	26.458	278	5994	0.353	ng/u1	96
29) Benzo(g,h,i)perylene	27.213	276	6796	0.366	ng/u1	97

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

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