

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012623\
 Data File : BM038574.D
 Acq On : 27 Jan 2023 07:39
 Operator : CG/JU
 Sample : PB150408BS
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS408

Quant Time: Jan 27 08:09:22 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	1375	0.400	ng/u1	0.00
4) Naphthalene-d8	10.774	136	3596	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.597	164	2156	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.341	188	3628	0.400	ng/u1	0.00
17) Chrysene-d12	21.513	240	2996	0.400	ng/u1	0.00
23) Perylene-d12	23.924	264	3788	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.343	96	1089	0.603	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.357	152	1790	0.321	ng/u1	0.00
18) Fluoranthene-d10	19.362	212	3540	0.414	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.377	88	2881	1.572	ng/u1#	20
5) Naphthalene	10.823	128	3016	0.258	ng/u1	99
7) 2-Methylnaphthalene	12.429	142	1920	0.311	ng/u1	100
8) 1-Methylnaphthalene	12.649	142	2053	0.322	ng/u1	99
10) Acenaphthylene	14.319	152	2952	0.310	ng/u1	99
11) Acenaphthene	14.657	153	2268	0.313	ng/u1	99
12) Fluorene	15.642	166	2615	0.289	ng/u1	99
14) Pentachlorophenol	16.995	266	801	0.542	ng/u1	99
15) Phenanthrene	17.379	178	3295	0.281	ng/u1	95
16) Anthracene	17.477	178	3222	0.322	ng/u1	97
19) Fluoranthene	19.394	202	4488	0.420	ng/u1	99
20) Pyrene	19.757	202	4692	0.408	ng/u1	97
21) Benzo(a)anthracene	21.498	228	3502	0.329	ng/u1	98
22) Chrysene	21.551	228	3646	0.341	ng/u1	99
24) Benzo(b)fluoranthene	23.184	252	4591	0.322	ng/u1	95
25) Benzo(k)fluoranthene	23.234	252	4528	0.327	ng/u1	96
26) Benzo(a)pyrene	23.816	252	4435	0.347	ng/u1	96
27) Indeno(1,2,3-cd)pyrene	26.431	276	6964	0.386	ng/u1#	65
28) Dibenzo(a,h)anthracene	26.445	278	5564	0.388	ng/u1	96
29) Benzo(g,h,i)perylene	27.203	276	6299	0.403	ng/u1	97

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

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