

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030923\
 Data File : BM038968.D
 Acq On : 10 Mar 2023 05:57
 Operator : CG/JU
 Sample : PB151192BS
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS192

Quant Time: Mar 10 06:29:07 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM030323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 09 13:36:03 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.995	152	10302	0.400	ng/u1	0.00
4) Naphthalene-d8	10.804	136	34052	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.627	164	19009	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.373	188	39522	0.400	ng/u1	0.00
17) Chrysene-d12	21.551	240	29872	0.400	ng/u1	0.00
23) Perylene-d12	23.995	264	20295	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	7391	0.634	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.388	152	14329	0.348	ng/u1	0.00
18) Fluoranthene-d10	19.397	212	28143	0.396	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.413	88	20438	1.732	ng/u1	93
5) Naphthalene	10.854	128	26630	0.327	ng/u1	100
7) 2-Methylnaphthalene	12.459	142	16628	0.337	ng/u1	100
8) 1-Methylnaphthalene	12.679	142	17303	0.334	ng/u1	99
10) Acenaphthylene	14.349	152	21200	0.307	ng/u1	99
11) Acenaphthene	14.691	153	18296	0.322	ng/u1	98
12) Fluorene	15.672	166	20658	0.322	ng/u1	99
14) Pentachlorophenol	17.018	266	4583	0.482	ng/u1	99
15) Phenanthrene	17.411	178	34190	0.327	ng/u1	99
16) Anthracene	17.504	178	28494	0.331	ng/u1	99
19) Fluoranthene	19.425	202	37977	0.375	ng/u1	98
20) Pyrene	19.787	202	38735	0.373	ng/u1	100
21) Benzo(a)anthracene	21.534	228	31430	0.372	ng/u1	100
22) Chrysene	21.589	228	34958	0.344	ng/u1	100
24) Benzo(b)fluoranthene	23.247	252	29808	0.376	ng/u1	98
25) Benzo(k)fluoranthene	23.296	252	28591	0.360	ng/u1	99
26) Benzo(a)pyrene	23.887	252	22042	0.349	ng/u1	96
27) Indeno(1,2,3-cd)pyrene	26.540	276	25115	0.314	ng/u1#	97
28) Dibenzo(a,h)anthracene	26.560	278	20942	0.344	ng/u1	98
29) Benzo(g,h,i)perylene	27.318	276	22508	0.302	ng/u1	99

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

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