

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030923\
 Data File : BM038969.D
 Acq On : 10 Mar 2023 06:34
 Operator : CG/JU
 Sample : PB151194BS
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS194

Quant Time: Mar 10 07:05:50 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	7.991	152	9011	0.400	ng/u1	0.00
4) Naphthalene-d8	10.805	136	29740	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.628	164	16541	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.370	188	34950	0.400	ng/u1	0.00
17) Chrysene-d12	21.553	240	29218	0.400	ng/u1	0.00
23) Perylene-d12	23.996	264	19177	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	7330	0.719	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.389	152	13914	0.387	ng/u1	0.00
18) Fluoranthene-d10	19.393	212	27187	0.391	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.413	88	18933	1.834	ng/u1	92
5) Naphthalene	10.855	128	25293	0.356	ng/u1	100
7) 2-Methylnaphthalene	12.461	142	15529	0.360	ng/u1	100
8) 1-Methylnaphthalene	12.681	142	16510	0.365	ng/u1	99
10) Acenaphthylene	14.350	152	20204	0.336	ng/u1	100
11) Acenaphthene	14.693	153	17367	0.351	ng/u1	98
12) Fluorene	15.673	166	19662	0.352	ng/u1	99
14) Pentachlorophenol	17.020	266	4445	0.529	ng/u1	98
15) Phenanthrene	17.412	178	32553	0.352	ng/u1	99
16) Anthracene	17.505	178	26966	0.354	ng/u1	99
19) Fluoranthene	19.426	202	36386	0.368	ng/u1	99
20) Pyrene	19.788	202	36752	0.362	ng/u1	98
21) Benzo(a)anthracene	21.535	228	29914	0.362	ng/u1	99
22) Chrysene	21.591	228	33656	0.339	ng/u1	100
24) Benzo(b)fluoranthene	23.248	252	28790	0.384	ng/u1	98
25) Benzo(k)fluoranthene	23.298	252	27108	0.362	ng/u1	98
26) Benzo(a)pyrene	23.888	252	20893	0.350	ng/u1	96
27) Indeno(1,2,3-cd)pyrene	26.537	276	24577	0.325	ng/u1#	97
28) Dibenzo(a,h)anthracene	26.557	278	20081	0.349	ng/u1	98
29) Benzo(g,h,i)perylene	27.319	276	21683	0.308	ng/u1	99

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

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