

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030923\
 Data File : BM038944.D
 Acq On : 09 Mar 2023 14:09
 Operator : CG/JU
 Sample : 01784-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCAD1

Quant Time: Mar 09 14:44:17 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	7.999	152	10229	0.400	ng/u1	0.00
4) Naphthalene-d8	10.810	136	35286	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.631	164	19769	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.373	188	41648	0.400	ng/u1	0.00
17) Chrysene-d12	21.556	240	32154	0.400	ng/u1	0.00
23) Perylene-d12	24.003	264	24738	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	55027	4.754	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.393	152	14439	0.338	ng/u1	0.00
18) Fluoranthene-d10	19.397	212	31811	0.416	ng/u1	0.00
Target Compounds						Qvalue
5) Naphthalene	10.859	128	12809	0.152	ng/u1	99
7) 2-Methylnaphthalene	12.465	142	1164	0.023	ng/u1	99
11) Acenaphthene	14.696	153	1360	0.023	ng/u1	100
14) Pentachlorophenol	17.023	266	737	0.074	ng/u1	98

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

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