

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032023\
 Data File : BM039055.D
 Acq On : 20 Mar 2023 21:48
 Operator : CG/JU
 Sample : 01964-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 UNION-2

Quant Time: Mar 21 01:51:44 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM030823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 17 05:25:34 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	316924	20.000	ng	0.00
21) Naphthalene-d8	10.786	136	1232258	20.000	ng	0.00
39) Acenaphthene-d10	14.610	164	691494	20.000	ng	0.00
64) Phenanthrene-d10	17.356	188	1451081	20.000	ng	0.00
76) Chrysene-d12	21.533	240	1391644	20.000	ng	0.00
86) Perylene-d12	23.968	264	1539838	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.528	112	3262856	156.417	ng	0.00
7) Phenol-d6	7.134	99	3797565	140.157	ng	0.00
23) Nitrobenzene-d5	9.139	82	2584556	103.027	ng	0.00
42) 2,4,6-Tribromophenol	16.098	330	1383037	172.891	ng	0.00
45) 2-Fluorobiphenyl	13.239	172	5216909	98.213	ng	0.00
79) Terphenyl-d14	19.980	244	8391811	110.251	ng	0.00

Target Compounds	Qvalue

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

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