

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032023\
 Data File : BM039058.D
 Acq On : 20 Mar 2023 23:42
 Operator : CG/JU
 Sample : 01964-11
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MAPLE-6

Quant Time: Mar 21 01:53:22 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	301466	20.000	ng	0.00
21) Naphthalene-d8	10.781	136	1149025	20.000	ng	-0.01
39) Acenaphthene-d10	14.610	164	620170	20.000	ng	0.00
64) Phenanthrene-d10	17.351	188	1199724	20.000	ng	-0.01
76) Chrysene-d12	21.533	240	1045287	20.000	ng	0.00
86) Perylene-d12	23.968	264	1187557	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.528	112	2950985	148.720	ng	0.00
7) Phenol-d6	7.134	99	3401474	131.976	ng	0.00
23) Nitrobenzene-d5	9.139	82	2360583	100.915	ng	0.00
42) 2,4,6-Tribromophenol	16.098	330	1053774	146.880	ng	0.00
45) 2-Fluorobiphenyl	13.239	172	4576872	96.073	ng	0.00
79) Terphenyl-d14	19.974	244	6325845	110.647	ng	0.00

Target Compounds	Qvalue

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

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