

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032023\
 Data File : BM039057.D
 Acq On : 20 Mar 2023 23:04
 Operator : CG/JU
 Sample : 01964-08
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MAPLE-5

Quant Time: Mar 21 01:52:33 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	286800	20.000	ng	0.00
21) Naphthalene-d8	10.781	136	1097669	20.000	ng	-0.01
39) Acenaphthene-d10	14.610	164	582665	20.000	ng	0.00
64) Phenanthrene-d10	17.351	188	1144181	20.000	ng	-0.01
76) Chrysene-d12	21.533	240	1044139	20.000	ng	0.00
86) Perylene-d12	23.962	264	1216268	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.528	112	2782462	147.398	ng	0.00
7) Phenol-d6	7.134	99	3144464	128.243	ng	0.00
23) Nitrobenzene-d5	9.140	82	2242786	100.365	ng	0.00
42) 2,4,6-Tribromophenol	16.098	330	1001464	148.574	ng	0.00
45) 2-Fluorobiphenyl	13.239	172	4342447	97.019	ng	0.00
79) Terphenyl-d14	19.974	244	6090113	106.640	ng	0.00
Target Compounds						
77) Benzdine	19.598	184	15598	5.465	ng	Qvalue 95

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

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