

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032223\
 Data File : BM039100.D
 Acq On : 22 Mar 2023 20:05
 Operator : CG/JU
 Sample : 02016-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 032023-1

Quant Time: Mar 23 04:26:05 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM030823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 22 03:36:20 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	257801	20.000	ng	0.00
21) Naphthalene-d8	10.781	136	957168	20.000	ng	0.00
39) Acenaphthene-d10	14.604	164	508481	20.000	ng	0.00
64) Phenanthrene-d10	17.351	188	995305	20.000	ng	0.00
76) Chrysene-d12	21.527	240	952752	20.000	ng	0.00
86) Perylene-d12	23.962	264	1042221	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	2492300	146.878	ng	0.00
7) Phenol-d6	7.128	99	2739801	124.308	ng	0.00
23) Nitrobenzene-d5	9.134	82	1995697	102.417	ng	0.00
42) 2,4,6-Tribromophenol	16.092	330	930613	158.205	ng	0.00
45) 2-Fluorobiphenyl	13.233	172	3852632	98.634	ng	0.00
79) Terphenyl-d14	19.968	244	5530693	106.134	ng	0.00
Target Compounds						
77) Benzdine	19.598	184	60183	7.366	ng	Qvalue 97

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

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