

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060923\
 Data File : BM040206.D
 Acq On : 10 Jun 2023 01:51
 Operator : MA/JU
 Sample : 03047-05
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB33B

Quant Time: Jun 10 04:46:19 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 08 23:25:50 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.987	152	282998	20.000	ng	0.00
21) Naphthalene-d8	10.798	136	1248252	20.000	ng	0.00
39) Acenaphthene-d10	14.621	164	792377	20.000	ng	0.00
64) Phenanthrene-d10	17.362	188	1814948	20.000	ng	0.00
76) Chrysene-d12	21.539	240	1720509	20.000	ng	-0.01
86) Perylene-d12	23.974	264	1845414	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.534	112	1318280	69.555	ng	0.00
7) Phenol-d6	7.140	99	1778578	68.772	ng	0.00
23) Nitrobenzene-d5	9.151	82	909556	39.439	ng	0.00
42) 2,4,6-Tribromophenol	16.110	330	843283	70.735	ng	0.00
45) 2-Fluorobiphenyl	13.251	172	2551134	42.554	ng	0.00
79) Terphenyl-d14	19.980	244	4991394	53.479	ng	0.00

Target Compounds	Qvalue

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

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