

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110123\
 Data File : BM042548.D
 Acq On : 01 Nov 2023 20:13
 Operator : MA/JU
 Sample : 05162-01 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 HD-01-10302023

Quant Time: Nov 01 23:44:57 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM103023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 31 02:55:18 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	73702	20.000	ng	0.00
21) Naphthalene-d8	10.745	136	273197	20.000	ng	0.00
39) Acenaphthene-d10	14.574	164	167676	20.000	ng	0.00
64) Phenanthrene-d10	17.327	188	362126	20.000	ng	# 0.00
76) Chrysene-d12	21.509	240	343217	20.000	ng	0.00
86) Perylene-d12	23.933	264	381273	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.457	112	191331	41.896	ng	0.00
7) Phenol-d6	7.087	99	246484	39.355	ng	-0.01
23) Nitrobenzene-d5	9.122	82	166331	28.382	ng	-0.01
42) 2,4,6-Tribromophenol	16.068	330	95321	43.542	ng	0.00
45) 2-Fluorobiphenyl	13.192	172	313440	25.154	ng	0.00
79) Terphenyl-d14	19.939	244	446642	22.066	ng	0.00
Target Compounds						
75) Fluoranthene	19.380	202	77433	3.483	ng	97
78) Pyrene	19.750	202	68073	2.888	ng	95

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

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