

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM052522\
 Data File : BM035247.D
 Acq On : 25 May 2022 18:33
 Operator : CG/JU
 Sample : N2922-09
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P024-SS001-2430-01

Quant Time: May 26 05:35:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 26 05:05:10 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	108626	20.000	ng	0.00
21) Naphthalene-d8	10.692	136	417333	20.000	ng	0.00
39) Acenaphthene-d10	14.527	164	265575	20.000	ng	0.00
64) Phenanthrene-d10	17.268	188	567918	20.000	ng	0.00
76) Chrysene-d12	21.450	240	595849	20.000	ng	# 0.00
86) Perylene-d12	23.827	264	633163	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	679723	110.395	ng	0.00
7) Phenol-d6	7.063	99	872315	109.745	ng	0.00
23) Nitrobenzene-d5	9.045	82	568728	73.592	ng	0.00
42) 2,4,6-Tribromophenol	16.015	330	371151	128.474	ng	0.00
45) 2-Fluorobiphenyl	13.151	172	1338133	73.144	ng	0.00
79) Terphenyl-d14	19.898	244	2253069	73.213	ng	0.00

Target Compounds	Qvalue

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

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