

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060822\
 Data File : BM035438.D
 Acq On : 08 Jun 2022 20:15
 Operator : CG/JU
 Sample : N3209-07
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-14B-10

Quant Time: Jun 09 00:51:24 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 07 14:28:07 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.839	152	171388	20.000	ng	0.00
21) Naphthalene-d8	10.639	136	615739	20.000	ng	# 0.00
39) Acenaphthene-d10	14.480	164	400208	20.000	ng	0.00
64) Phenanthrene-d10	17.227	188	812505	20.000	ng	0.00
76) Chrysene-d12	21.415	240	744970	20.000	ng	# 0.00
86) Perylene-d12	23.768	264	854475	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	1245060	152.652	ng	0.00
7) Phenol-d6	7.022	99	1355701	121.685	ng	0.00
23) Nitrobenzene-d5	9.004	82	997996	102.915	ng	0.00
42) 2,4,6-Tribromophenol	15.974	330	853427	132.907	ng	0.00
45) 2-Fluorobiphenyl	13.104	172	2937757	105.430	ng	0.00
79) Terphenyl-d14	19.850	244	4197841	112.263	ng	0.00

Target Compounds	Qvalue

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

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