

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080123\
 Data File : BP016455.D
 Acq On : 01 Aug 2023 14:29
 Operator : MA/JU
 Sample : 03810-09
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 B-P1B

Quant Time: Aug 01 16:32:09 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 28 01:48:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.081	152	510571	20.000	ng	-0.01
21) Naphthalene-d8	10.910	136	2110413	20.000	ng	-0.02
39) Acenaphthene-d10	14.728	164	1248976	20.000	ng	0.00
64) Phenanthrene-d10	17.492	188	2667209	20.000	ng	0.00
76) Chrysene-d12	21.586	240	1847617	20.000	ng	# 0.00
86) Perylene-d12	24.186	264	1747363	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.587	112	3221664	102.941	ng	-0.02
7) Phenol-d6	7.216	99	4305199	100.361	ng	-0.01
23) Nitrobenzene-d5	9.275	82	2444218	64.833	ng	-0.02
42) 2,4,6-Tribromophenol	16.222	330	1346475	73.281	ng	0.00
45) 2-Fluorobiphenyl	13.351	172	5036320	57.811	ng	-0.01
79) Terphenyl-d14	20.039	244	7405732	84.671	ng	0.00

Target Compounds	Qvalue

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

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