

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080823\
 Data File : BP016538.D
 Acq On : 08 Aug 2023 14:59
 Operator : MA/JU
 Sample : 03895-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MCLEAN-TP11

Quant Time: Aug 08 16:07:39 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 08 16:03:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.063	152	214290	20.000	ng	0.00
21) Naphthalene-d8	10.898	136	903597	20.000	ng	0.00
39) Acenaphthene-d10	14.716	164	581962	20.000	ng	0.00
64) Phenanthrene-d10	17.480	188	1499126	20.000	ng	0.00
76) Chrysene-d12	21.574	240	1681489	20.000	ng	# 0.00
86) Perylene-d12	24.168	264	1894081	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	2382288	181.366	ng	0.00
7) Phenol-d6	7.204	99	3051709	169.500	ng	0.00
23) Nitrobenzene-d5	9.263	82	1936126	119.944	ng	0.00
42) 2,4,6-Tribromophenol	16.210	330	1383252	161.566	ng	0.00
45) 2-Fluorobiphenyl	13.339	172	4450247	109.632	ng	0.00
79) Terphenyl-d14	20.027	244	8998993	113.052	ng	0.00

Target Compounds	Qvalue

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

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