

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061021\
 Data File : BP005863.D
 Acq On : 11 Jun 2021 09:15
 Operator : CG/JU
 Sample : M2612-39
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B12(78-82)

Quant Time: Jun 11 11:04:58 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 09 11:54:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	133009	20.000	ng	-0.01
21) Naphthalene-d8	10.751	136	469764	20.000	ng	-0.01
39) Acenaphthene-d10	14.575	164	271493	20.000	ng	0.00
64) Phenanthrene-d10	17.322	188	572430	20.000	ng	0.00
76) Chrysene-d12	21.392	240	699618	20.000	ng	0.00
86) Perylene-d12	23.821	264	678233	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	1117158	134.775	ng	0.00
7) Phenol-d6	7.116	99	1170010	108.900	ng	0.00
23) Nitrobenzene-d5	9.110	82	875250	91.346	ng	-0.01
42) 2,4,6-Tribromophenol	16.063	330	661102	141.897	ng	0.00
45) 2-Fluorobiphenyl	13.204	172	1827693	88.400	ng	-0.01
79) Terphenyl-d14	19.874	244	3271115	87.550	ng	0.00

Target Compounds	Qvalue

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

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