

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061021\
 Data File : BP005857.D
 Acq On : 11 Jun 2021 04:22
 Operator : CG/JU
 Sample : M2612-29
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B12(38-42)

Quant Time: Jun 11 07:03:42 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	70836	20.000	ng	-0.01
21) Naphthalene-d8	10.751	136	256434	20.000	ng	-0.01
39) Acenaphthene-d10	14.575	164	153352	20.000	ng	0.00
64) Phenanthrene-d10	17.316	188	321894	20.000	ng	-0.01
76) Chrysene-d12	21.386	240	365382	20.000	ng	0.00
86) Perylene-d12	23.804	264	440442	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.511	112	568641	128.813	ng	0.00
7) Phenol-d6	7.110	99	573750	100.274	ng	-0.01
23) Nitrobenzene-d5	9.110	82	474172	90.656	ng	-0.01
42) 2,4,6-Tribromophenol	16.063	330	355348	135.029	ng	0.00
45) 2-Fluorobiphenyl	13.204	172	1024297	87.709	ng	-0.01
79) Terphenyl-d14	19.863	244	1732874	88.805	ng	0.00

Target Compounds	Qvalue

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

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