

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061021\
 Data File : BP005861.D
 Acq On : 11 Jun 2021 06:37
 Operator : CG/JU
 Sample : M2612-35
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B12(62-66)

Quant Time: Jun 11 08:45:33 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	7.945	152	69261	20.000	ng	-0.01
21) Naphthalene-d8	10.751	136	251453	20.000	ng	-0.01
39) Acenaphthene-d10	14.575	164	149560	20.000	ng	0.00
64) Phenanthrene-d10	17.316	188	320740	20.000	ng	-0.01
76) Chrysene-d12	21.386	240	373055	20.000	ng	0.00
86) Perylene-d12	23.803	264	448745	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	555278	128.646	ng	0.00
7) Phenol-d6	7.110	99	580578	103.775	ng	-0.01
23) Nitrobenzene-d5	9.110	82	452099	88.148	ng	-0.01
42) 2,4,6-Tribromophenol	16.057	330	339026	132.093	ng	-0.01
45) 2-Fluorobiphenyl	13.204	172	969751	85.144	ng	-0.01
79) Terphenyl-d14	19.862	244	1711267	85.894	ng	0.00

Target Compounds	Qvalue

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

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