

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060821\
 Data File : BP005793.D
 Acq On : 09 Jun 2021 03:37
 Operator : CG/JU
 Sample : M2589-36
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B11(4-6)

Quant Time: Jun 09 06:01:08 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 02:47:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	72262	20.000	ng	0.00
21) Naphthalene-d8	10.763	136	256515	20.000	ng	0.00
39) Acenaphthene-d10	14.581	164	151638	20.000	ng	0.00
64) Phenanthrene-d10	17.322	188	307530	20.000	ng	0.00
76) Chrysene-d12	21.392	240	352318	20.000	ng	0.00
86) Perylene-d12	23.815	264	421382	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	627829	139.414	ng	0.00
7) Phenol-d6	7.122	99	666831	114.242	ng	0.00
23) Nitrobenzene-d5	9.122	82	491446	93.929	ng	0.00
42) 2,4,6-Tribromophenol	16.069	330	348982	134.109	ng	0.00
45) 2-Fluorobiphenyl	13.210	172	1060970	91.876	ng	0.00
79) Terphenyl-d14	19.869	244	1675652	89.057	ng	0.00

Target Compounds	Qvalue

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

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