

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060921\
 Data File : BP005816.D
 Acq On : 09 Jun 2021 19:10
 Operator : CG/JU
 Sample : M2589-35
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B11(2-4)

Quant Time: Jun 10 01:21:46 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	7.957	152	69963	20.000	ng	0.00
21) Naphthalene-d8	10.757	136	256607	20.000	ng	0.00
39) Acenaphthene-d10	14.581	164	155862	20.000	ng	0.00
64) Phenanthrene-d10	17.322	188	324196	20.000	ng	0.00
76) Chrysene-d12	21.392	240	394400	20.000	ng	0.00
86) Perylene-d12	23.815	264	475047	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	568634	130.418	ng	0.00
7) Phenol-d6	7.116	99	609154	107.790	ng	0.00
23) Nitrobenzene-d5	9.116	82	451152	86.197	ng	0.00
42) 2,4,6-Tribromophenol	16.063	330	362368	135.479	ng	0.00
45) 2-Fluorobiphenyl	13.210	172	970065	81.728	ng	0.00
79) Terphenyl-d14	19.869	244	1754047	83.277	ng	0.00

Target Compounds	Qvalue

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

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