

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061621\
 Data File : BP005953.D
 Acq On : 16 Jun 2021 17:41
 Operator : CG/JU
 Sample : M2672-32
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 17-B6(52-56)

Quant Time: Jun 17 01:57:33 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 16 11:35:30 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	64186	20.000	ng	0.00
21) Naphthalene-d8	10.734	136	229228	20.000	ng	0.00
39) Acenaphthene-d10	14.557	164	134035	20.000	ng	0.00
64) Phenanthrene-d10	17.304	188	291189	20.000	ng	0.00
76) Chrysene-d12	21.374	240	372663	20.000	ng	0.00
86) Perylene-d12	23.786	264	432277	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	532160	133.038	ng	0.00
7) Phenol-d6	7.104	99	567107	109.382	ng	0.00
23) Nitrobenzene-d5	9.093	82	412394	88.203	ng	0.00
42) 2,4,6-Tribromophenol	16.045	330	330673	143.762	ng	0.00
45) 2-Fluorobiphenyl	13.186	172	915314	89.673	ng	0.00
79) Terphenyl-d14	19.851	244	1798276	90.356	ng	0.00

Target Compounds	Qvalue

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

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