

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061621\
 Data File : BP005957.D
 Acq On : 16 Jun 2021 19:57
 Operator : CG/JU
 Sample : M2672-22
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 17-B6(12-16)

Quant Time: Jun 17 01:58:50 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	68532	20.000	ng	0.00
21) Naphthalene-d8	10.734	136	247528	20.000	ng	0.00
39) Acenaphthene-d10	14.557	164	150915	20.000	ng	0.00
64) Phenanthrene-d10	17.304	188	324246	20.000	ng	0.00
76) Chrysene-d12	21.374	240	405900	20.000	ng	0.00
86) Perylene-d12	23.786	264	459269	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	556221	130.235	ng	0.00
7) Phenol-d6	7.104	99	582521	105.230	ng	0.00
23) Nitrobenzene-d5	9.092	82	441528	87.452	ng	0.00
42) 2,4,6-Tribromophenol	16.045	330	370230	142.956	ng	0.00
45) 2-Fluorobiphenyl	13.186	172	1008963	87.791	ng	0.00
79) Terphenyl-d14	19.851	244	1939067	89.453	ng	0.00

Target Compounds	Qvalue

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

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