

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061021\
 Data File : BP005847.D
 Acq On : 10 Jun 2021 20:26
 Operator : CG/JU
 Sample : M2657-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 HT-3-PED

Quant Time: Jun 11 07:01:28 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	7.952	152	73495	20.000	ng	0.00
21) Naphthalene-d8	10.757	136	271113	20.000	ng	0.00
39) Acenaphthene-d10	14.575	164	172852	20.000	ng	0.00
64) Phenanthrene-d10	17.322	188	380013	20.000	ng	0.00
76) Chrysene-d12	21.386	240	453806	20.000	ng	0.00
86) Perylene-d12	23.810	264	517254	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.511	112	338359	73.874	ng	0.00
7) Phenol-d6	7.110	99	520268	87.638	ng	-0.01
23) Nitrobenzene-d5	9.110	82	341756	61.802	ng	-0.01
42) 2,4,6-Tribromophenol	16.063	330	184311	62.136	ng	0.00
45) 2-Fluorobiphenyl	13.204	172	817277	62.087	ng	-0.01
79) Terphenyl-d14	19.869	244	1517305	62.607	ng	0.00

Target Compounds	Qvalue

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

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