

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061421\
 Data File : BP005905.D
 Acq On : 14 Jun 2021 14:39
 Operator : CG/JU
 Sample : M2637-07
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B12(234-238)

Quant Time: Jun 14 15:11:02 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 14 13:15:03 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	71774	20.000	ng	0.00
21) Naphthalene-d8	10.740	136	264233	20.000	ng	0.00
39) Acenaphthene-d10	14.563	164	161323	20.000	ng	0.00
64) Phenanthrene-d10	17.310	188	345526	20.000	ng	0.00
76) Chrysene-d12	21.380	240	419181	20.000	ng	0.00
86) Perylene-d12	23.792	264	480477	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.511	112	616176	137.756	ng	0.00
7) Phenol-d6	7.116	99	653514	112.722	ng	0.00
23) Nitrobenzene-d5	9.099	82	491708	91.234	ng	0.00
42) 2,4,6-Tribromophenol	16.051	330	417752	150.899	ng	0.00
45) 2-Fluorobiphenyl	13.193	172	1094970	89.128	ng	0.00
79) Terphenyl-d14	19.857	244	2054476	91.774	ng	0.00
Target Compounds						
32) Benzoic acid	10.040	122	3211	7.255	ng	97
84) Bis(2-ethylhexyl)phtha...	21.286	149	106734	6.045	ng	99

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

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