

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061021\
 Data File : BP005837.D
 Acq On : 10 Jun 2021 14:45
 Operator : CG/JU
 Sample : M2613-25
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B12(130-134)

Manual Integrations
 APPROVED

mohammad
 6/11/2021 2:12:30 PM

Quant Time: Jun 10 15:43:18 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	66442	20.000	ng	0.00
21) Naphthalene-d8	10.757	136	234941	20.000	ng	0.00
39) Acenaphthene-d10	14.575	164	142102	20.000	ng	0.00
64) Phenanthrene-d10	17.322	188	299262	20.000	ng	0.00
76) Chrysene-d12	21.392	240	358127	20.000	ng	0.00
86) Perylene-d12	23.810	264	431208	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	497281	120.097	ng	0.00
7) Phenol-d6	7.116	99	533327	99.374	ng	0.00
23) Nitrobenzene-d5	9.116	82	384645	80.267	ng	0.00
42) 2,4,6-Tribromophenol	16.063	330	308111	126.349	ng	0.00
45) 2-Fluorobiphenyl	13.204	172	853554	78.875	ng	-0.01
79) Terphenyl-d14	19.869	244	1494463	78.139	ng	0.00
Target Compounds						Qvalue
32) Benzoic acid	10.087	122	1067m	6.669	ng	

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

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