

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060821\
 Data File : BP005808.D
 Acq On : 09 Jun 2021 12:08
 Operator : CG/JU
 Sample : M2589-33
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B10(4-6)

Manual Integrations
 APPROVED

mohammad
 6/9/2021 3:53:31 PM

Quant Time: Jun 09 12:54:10 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 02:47:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	74536	20.000	ng	0.00
21) Naphthalene-d8	10.763	136	271117	20.000	ng	0.00
39) Acenaphthene-d10	14.581	164	161839	20.000	ng	0.00
64) Phenanthrene-d10	17.328	188	341493	20.000	ng	0.00
76) Chrysene-d12	21.398	240	398653	20.000	ng	0.00
86) Perylene-d12	23.821	264	476143	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	629904	135.607	ng	0.00
7) Phenol-d6	7.122	99	661524	109.875	ng	0.00
23) Nitrobenzene-d5	9.122	82	515186	93.163	ng	0.00
42) 2,4,6-Tribromophenol	16.069	330	386981	139.338	ng	0.00
45) 2-Fluorobiphenyl	13.210	172	1107151	89.832	ng	0.00
79) Terphenyl-d14	19.875	244	1868912	87.784	ng	0.00
Target Compounds						
31) Naphthalene	10.810	128	46125	2.940	ng	Qvalue 98
32) Benzoic acid	10.099	122	2181m	6.939	ng	

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

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