

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060821\
 Data File : BP005797.D
 Acq On : 09 Jun 2021 05:53
 Operator : CG/JU
 Sample : M2589-22
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B7(2-4)

Manual Integrations
 APPROVED

mohammad
 6/9/2021 3:53:27 PM

Quant Time: Jun 09 07:25:55 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	79110	20.000	ng	0.00
21) Naphthalene-d8	10.763	136	284445	20.000	ng	0.00
39) Acenaphthene-d10	14.581	164	171819	20.000	ng	0.00
64) Phenanthrene-d10	17.322	188	361673	20.000	ng	0.00
76) Chrysene-d12	21.392	240	407459	20.000	ng	0.00
86) Perylene-d12	23.816	264	481612	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	648128	131.463	ng	0.00
7) Phenol-d6	7.122	99	685449	107.267	ng	0.00
23) Nitrobenzene-d5	9.122	82	519866	89.604	ng	0.00
42) 2,4,6-Tribromophenol	16.069	330	398169	135.039	ng	0.00
45) 2-Fluorobiphenyl	13.210	172	1140339	87.151	ng	0.00
79) Terphenyl-d14	19.869	244	1901130	87.367	ng	0.00
Target Compounds						Qvalue
32) Benzoic acid	10.069	122	2552m	7.010	ng	

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

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