

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052922\
 Data File : BP010618.D
 Acq On : 29 May 2022 14:39
 Operator : CG/JU
 Sample : N2859-09
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SW-E-E3

Manual Integrations
 APPROVED

Reviewed By :Christian
 Giraldo

Quant Time: May 30 02:06:15 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 28 02:18:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	134590	20.000	ng	0.00
21) Naphthalene-d8	10.669	136	525521	20.000	ng	0.00
39) Acenaphthene-d10	14.498	164	333636	20.000	ng	0.00
64) Phenanthrene-d10	17.257	188	704612	20.000	ng	0.00
76) Chrysene-d12	21.339	240	666358	20.000	ng	0.00
86) Perylene-d12	23.751	264	669689	20.000	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.452	112	461003	62.487	ng	0.00
7) Phenol-d6	7.040	99	356109	34.001	ng	0.00
23) Nitrobenzene-d5	9.028	82	1011964	91.041	ng	0.00
42) 2,4,6-Tribromophenol	15.998	330	555548	147.102	ng	0.00
45) 2-Fluorobiphenyl	13.128	172	2056440	87.478	ng	0.00
79) Terphenyl-d14	19.816	244	3311459	93.342	ng	0.00

Target Compounds						Qvalue
32) Benzoic acid	9.975	122	3363m	7.920	ng	

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

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