

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052922\
 Data File : BP010627.D
 Acq On : 29 May 2022 19:44
 Operator : CG/JU
 Sample : N3056-09DL 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-02DL

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 05/31/2022
 Supervised By :Jagrut Upadhyay 05/31/2022

Quant Time: May 30 02:07:42 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	177483	20.000	ng	0.00
21) Naphthalene-d8	10.669	136	683664	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	408015	20.000	ng	0.00
64) Phenanthrene-d10	17.257	188	758701	20.000	ng	0.00
76) Chrysene-d12	21.339	240	685872	20.000	ng	0.00
86) Perylene-d12	23.757	264	715518	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.452	112	127040	13.058	ng	0.00
7) Phenol-d6	7.052	99	100839	7.301	ng	0.01
23) Nitrobenzene-d5	9.028	82	225745	15.611	ng	0.00
42) 2,4,6-Tribromophenol	15.998	330	125287	27.127	ng	0.00
45) 2-Fluorobiphenyl	13.128	172	497950	17.321	ng	0.00
79) Terphenyl-d14	19.816	244	591228	16.191	ng	0.00
Target Compounds						Qvalue
27) 2,4-Dimethylphenol	9.858	122	27748m	3.279	ng	
31) Naphthalene	10.722	128	1547604	42.970	ng	99
37) 2-Methylnaphthalene	12.328	142	210360	8.371	ng	99
38) 1-Methylnaphthalene	12.546	142	181803m	7.408	ng	

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

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