

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052922\
 Data File : BP010631.D
 Acq On : 29 May 2022 21:59
 Operator : CG/JU
 Sample : N3056-02DL 20X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-12DL

Quant Time: May 30 02:08:23 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	171127	20.000	ng	0.00
21) Naphthalene-d8	10.669	136	680486	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	425771	20.000	ng	0.00
64) Phenanthrene-d10	17.257	188	889292	20.000	ng	0.00
76) Chrysene-d12	21.345	240	772460	20.000	ng	0.00
86) Perylene-d12	23.757	264	771833	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.452	112	27348	2.915	ng	0.00
7) Phenol-d6	7.052	99	19407	1.457	ng	0.01
23) Nitrobenzene-d5	9.034	82	44814	3.114	ng	0.00
42) 2,4,6-Tribromophenol	15.998	330	30480	6.324	ng	0.00
45) 2-Fluorobiphenyl	13.128	172	117980	3.933	ng	0.00
79) Terphenyl-d14	19.810	244	149952	3.646	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	10.716	128	1438578	40.130	ng	99
37) 2-Methylnaphthalene	12.328	142	157014	6.277	ng	98
38) 1-Methylnaphthalene	12.545	142	452084	18.507	ng	100
52) Acenaphthene	14.569	154	79468	3.376	ng	99

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

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