

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070820\
 Data File : BP002672.D
 Acq On : 08 Jul 2020 13:28
 Operator : CG/JU
 Sample : L3215-11DL 10X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-14DL

Quant Time: Jul 08 14:12:21 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 29 16:51:57 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	227487	20.00	ng	0.00
21) Naphthalene-d8	10.34	136	946524	20.00	ng	0.00
39) Acenaphthene-d10	14.22	164	607743	20.00	ng	0.00
64) Phenanthrene-d10	16.97	188	1403173	20.00	ng	0.00
76) Chrysene-d12	21.09	240	1638363	20.00	ng	0.00
86) Perylene-d12	23.28	264	1582007	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	97107	7.23	ng	0.00
7) Phenol-d6	6.77	99	78630	4.19	ng	0.00
23) Nitrobenzene-d5	8.72	82	186854	10.17	ng	0.00
42) 2,4,6-Tribromophenol	15.72	330	148009	20.02	ng	0.00
45) 2-Fluorobiphenyl	12.83	172	448609	10.86	ng	0.00
79) Terphenyl-d14	19.56	244	911949	12.01	ng	-0.01
Target Compounds						
31) Naphthalene	10.39	128	6772452	135.161	ng	98
37) 2-Methylnaphthalene	12.01	142	1065138	29.154	ng	97
38) 1-Methylnaphthalene	12.23	142	1660949	48.272	ng	98
46) 1,1'-Biphenyl	13.05	154	155858	3.390	ng	99
52) Acenaphthene	14.28	154	449376	12.933	ng	97
58) Fluorene	15.27	166	179388	4.183	ng	98
71) Phenanthrene	17.02	178	478066	6.332	ng	99

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

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