

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071620\
 Data File : BP002779.D
 Acq On : 16 Jul 2020 19:03
 Operator : CG/JU
 Sample : L3308-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 5

Quant Time: Jul 17 00:54:23 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 13 10:54:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	212752	20.00	ng	0.00
21) Naphthalene-d8	10.33	136	898709	20.00	ng	0.00
39) Acenaphthene-d10	14.21	164	530360	20.00	ng	0.00
64) Phenanthrene-d10	16.97	188	1018021	20.00	ng	0.00
76) Chrysene-d12	21.07	240	709386	20.00	ng	0.00
86) Perylene-d12	23.27	264	633936	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	1048256	83.13	ng	0.00
7) Phenol-d6	6.76	99	1443775	80.23	ng	0.00
23) Nitrobenzene-d5	8.70	82	930670	55.39	ng	0.00
42) 2,4,6-Tribromophenol	15.72	330	443664	80.14	ng	0.00
45) 2-Fluorobiphenyl	12.82	172	1746218	50.93	ng	0.00
79) Terphenyl-d14	19.56	244	1864921	60.27	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.67	163	353512	8.740	ng	99
71) Phenanthrene	17.01	178	121953	2.213	ng	98
75) Fluoranthene	19.00	202	314870	4.930	ng	98
78) Pyrene	19.35	202	234012	4.778	ng	98
83) Chrysene	21.12	228	112665	2.577	ng	97
88) Benzo(b)fluoranthene	22.62	252	132644	3.370	ng	# 94

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

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