

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
 Data File : BP001412.D
 Acq On : 25 Dec 2019 07:48
 Operator : JU
 Sample : K6372-18DL 10X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-14DL

Quant Time: Jan 01 06:43:43 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 16:04:49 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.29	152	34618	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	123385	20.00	ng	0.00
39) Acenaphthene-d10	13.18	164	68859	20.00	ng	0.00
64) Phenanthrene-d10	15.58	188	129841	20.00	ng	0.00
76) Chrysene-d12	19.22	240	109804	20.00	ng	0.00
87) Perylene-d12	20.99	264	110181	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.20	112	10395	4.85	ng	0.00
7) Phenol-d6	7.75	99	9323	3.34	ng	0.00
23) Nitrobenzene-d5	9.16	82	26403	8.22	ng	-0.01
42) 2,4,6-Tribromophenol	14.47	330	10031	11.65	ng	0.00
45) 2-Fluorobiphenyl	12.10	172	50039	9.19	ng	0.00
79) Terphenyl-d14	17.86	244	61942	9.66	ng	0.00
Target Compounds						
31) Naphthalene	10.37	128	1588383	235.833	ng	99
37) 2-Methylnaphthalene	11.49	142	154096	33.141	ng	99
38) 1-Methylnaphthalene	11.65	142	153707	35.162	ng	98
52) Acenaphthene	13.23	154	30565	7.450	ng	99
71) Phenanthrene	15.62	178	23441	3.164	ng	97

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

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