

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
 Data File : BP001393.D
 Acq On : 24 Dec 2019 18:35
 Operator : JU
 Sample : K6372-12
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-3

Quant Time: Jan 01 06:42:08 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	51238	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	189198	20.00	ng	0.00
39) Acenaphthene-d10	13.18	164	109051	20.00	ng	0.00
64) Phenanthrene-d10	15.58	188	202396	20.00	ng	0.00
76) Chrysene-d12	19.22	240	144610	20.00	ng	0.00
87) Perylene-d12	20.99	264	142198	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.20	112	164828	51.99	ng	0.00
7) Phenol-d6	7.75	99	125331	30.30	ng	0.00
23) Nitrobenzene-d5	9.18	82	423806	86.02	ng	0.00
42) 2,4,6-Tribromophenol	14.48	330	192786	141.40	ng	0.00
45) 2-Fluorobiphenyl	12.10	172	789862	91.59	ng	0.00
79) Terphenyl-d14	17.87	244	868560	102.87	ng	0.00
Target Compounds						
10) Phenol	7.77	94	11182	2.366	ng	91
22) Acetophenone	8.93	105	118251	19.765	ng	# 62
31) Naphthalene	10.36	128	571289	55.316	ng	99
32) Benzoic acid	9.96	122	4866	7.142	ng	92
37) 2-Methylnaphthalene	11.48	142	38555	5.407	ng	# 92
38) 1-Methylnaphthalene	11.65	142	103459	15.434	ng	# 98

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

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