

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021319\
 Data File : BM018811.D
 Acq On : 13 Feb 2019 20:48
 Operator : JU/SJ
 Sample : K1457-07
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 IDW-LIQUID-01

Manual Integrations
 APPROVED

Sohil
 2/14/2019 1:50:33 PM

Quant Time: Feb 14 03:40:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 11 16:02:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	299366	20.00	ng	0.00
21) Naphthalene-d8	10.50	136	1254909	20.00	ng	0.00
39) Acenaphthene-d10	14.36	164	705512	20.00	ng	0.00
64) Phenanthrene-d10	17.11	188	1610668	20.00	ng	0.00
76) Chrysene-d12	21.30	240	1693744	20.00	ng	0.00
87) Perylene-d12	23.56	264	1600092	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	2522018	138.03	ng	0.00
7) Phenol-d6	6.89	99	3129456	121.58	ng	0.00
23) Nitrobenzene-d5	8.88	82	2550256	93.97	ng	0.00
42) 2,4,6-Tribromophenol	15.86	330	1480895	145.62	ng	0.00
45) 2-Fluorobiphenyl	12.97	172	4860579	91.46	ng	0.00
79) Terphenyl-d14	19.74	244	8431210	102.69	ng	0.00
Target Compounds						
20) 3+4-Methylphenols	8.48	107	65607	2.756	ng	93
31) Naphthalene	10.55	128	891253	14.562	ng	99
32) Benzoic acid	9.77	122	83843m	5.866	ng	
37) 2-Methylnaphthalene	12.16	142	172962	4.014	ng	99

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

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