

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020719\
 Data File : BF112438.D
 Acq On : 7 Feb 2019 15:18
 Operator : JU/SJ
 Sample : K1389-11
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 M-5S

Manual Integrations
 APPROVED

Sohil
 2/11/2019 4:49:58 PM

Quant Time: Feb 09 03:32:27 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 28 10:38:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	155076	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	605507	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	286626	20.00	ng	-0.02
64) Phenanthrene-d10	11.36	188	503884	20.00	ng	-0.01
76) Chrysene-d12	14.00	240	356957	20.00	ng	-0.01
87) Perylene-d12	15.47	264	324140	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	921660	126.24	ng	0.00
7) Phenol-d6	6.49	99	1161959	114.54	ng	0.00
23) Nitrobenzene-d5	7.40	82	726782	69.16	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	333013	99.82	ng	-0.02
45) 2-Fluorobiphenyl	9.19	172	1324429	65.35	ng	-0.02
79) Terphenyl-d14	12.94	244	1141143	60.21	ng	-0.01
Target Compounds						
10) Phenol	6.49	94	23365	2.099	ng	Qvalue # 84
50) Dimethylphthalate	9.58	163	120754	5.422	ng	98
75) Fluoranthene	12.57	202	65807	2.342	ng	98
78) Pyrene	12.80	202	61324	2.481	ng	97
88) Benzo(b)fluoranthene	15.04	252	43160m	2.226	ng	

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

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