

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022719\
 Data File : BF112720.D
 Acq On : 27 Feb 2019 15:42
 Operator : JU/SJ
 Sample : K1613-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 R-1

Manual Integrations
 APPROVED

Sohil
 2/28/2019 12:46:58 PM

Quant Time: Feb 28 02:48:13 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 27 16:08:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	225732	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	829014	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	371344	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	656106	20.00	ng	0.00
76) Chrysene-d12	13.87	240	450792	20.00	ng	0.00
87) Perylene-d12	15.27	264	526306	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	1693724	116.72	ng	0.01
7) Phenol-d6	6.37	99	2182682	119.74	ng	0.00
23) Nitrobenzene-d5	7.30	82	1309652	94.53	ng	0.00
42) 2,4,6-Tribromophenol	10.55	330	431302	109.40	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	1937265	90.36	ng	0.00
79) Terphenyl-d14	12.83	244	1601205	68.86	ng	0.00
Target Compounds						
10) Phenol	6.39	94	74368	3.496	ng	88
50) Dimethylphthalate	9.49	163	157259	5.743	ng	98
78) Pyrene	12.67	202	90988	2.394	ng	98
84) Bis(2-ethylhexyl)phthalate	13.87	149	183467	9.325	ng	98
88) Benzo(b)fluoranthene	14.87	252	78378m	2.302	ng	

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

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