

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012419\
 Data File : BF112211.D
 Acq On : 24 Jan 2019 16:19
 Operator : JU/SJ
 Sample : K1016-07 2X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-05-RO-3-012319

Manual Integrations
 APPROVED

Sohil
 1/25/2019 10:05:05 AM

Quant Time: Jan 25 01:37:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 19 01:02:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	211606	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	810667	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	368938	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	595353	20.00	ng	0.00
76) Chrysene-d12	14.03	240	501089	20.00	ng	0.00
87) Perylene-d12	15.50	264	386053	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	633449	54.33	ng	0.00
7) Phenol-d6	6.50	99	795088	54.51	ng	0.00
23) Nitrobenzene-d5	7.41	82	483617	41.22	ng	-0.02
42) 2,4,6-Tribromophenol	10.68	330	197468	49.61	ng	-0.01
45) 2-Fluorobiphenyl	9.21	172	927803	36.87	ng	0.00
79) Terphenyl-d14	12.98	244	669013	28.40	ng	0.01
Target Compounds						
10) Phenol	6.51	94	68585	4.302	ng	Qvalue # 72
50) Dimethylphthalate	9.60	163	81912	3.044	ng	100
78) Pyrene	12.87	202	66190m	2.002	ng	
88) Benzo(b)fluoranthene	15.07	252	55616m	2.538	ng	

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

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