

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030619\
 Data File : BF112911.D
 Acq On : 7 Mar 2019 5:16
 Operator : JU/SJ
 Sample : K1705-18
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PST-028-B004-022719

Manual Integrations
 APPROVED

Sohil
 3/7/2019 11:03:07 AM

Quant Time: Mar 07 05:56:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 01 15:40:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	197138	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	701761	20.00	ng	-0.01
39) Acenaphthene-d10	9.76	164	290317	20.00	ng	-0.01
64) Phenanthrene-d10	11.24	188	511335	20.00	ng	0.00
76) Chrysene-d12	13.86	240	537544	20.00	ng	-0.01
87) Perylene-d12	15.26	264	486783	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1434262	113.17	ng	0.00
7) Phenol-d6	6.37	99	1784042	112.07	ng	0.00
23) Nitrobenzene-d5	7.29	82	1075492	91.71	ng	-0.01
42) 2,4,6-Tribromophenol	10.55	330	396156	128.54	ng	-0.01
45) 2-Fluorobiphenyl	9.09	172	1541672	89.13	ng	-0.01
79) Terphenyl-d14	12.82	244	1357829	48.97	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.48	163	123769	5.782	ng	98
71) Phenanthrene	11.26	178	69968	2.750	ng	96
75) Fluoranthene	12.45	202	111931	4.107	ng	96
78) Pyrene	12.67	202	110138	2.430	ng	98
88) Benzo(b)fluoranthene	14.87	252	69773m	2.216	ng	

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

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