

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030619\
 Data File : BF112910.D
 Acq On : 7 Mar 2019 4:49
 Operator : JU/SJ
 Sample : K1705-16
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PST-028-B101-022719

Manual Integrations
 APPROVED

Sohil
 3/7/2019 11:03:04 AM

Quant Time: Mar 07 05:53:41 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	193470	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	709862	20.00	ng	-0.01
39) Acenaphthene-d10	9.76	164	310070	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	576028	20.00	ng	0.00
76) Chrysene-d12	13.86	240	589886	20.00	ng	-0.01
87) Perylene-d12	15.27	264	531498	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1068064	85.87	ng	0.00
7) Phenol-d6	6.37	99	1422333	91.04	ng	0.00
23) Nitrobenzene-d5	7.29	82	808561	68.16	ng	-0.02
42) 2,4,6-Tribromophenol	10.55	330	354611	107.73	ng	-0.01
45) 2-Fluorobiphenyl	9.09	172	1372661	70.02	ng	-0.01
79) Terphenyl-d14	12.82	244	1267086	41.64	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.48	163	115286	5.042	ng	100
71) Phenanthrene	11.26	178	99920	3.486	ng	98
75) Fluoranthene	12.44	202	142434	4.639	ng	94
78) Pyrene	12.67	202	136127	2.737	ng	98
88) Benzo(b)fluoranthene	14.87	252	79076m	2.300	ng	

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

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