

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031219\
 Data File : BF113028.D
 Acq On : 13 Mar 2019 9:50
 Operator : JU/SJ
 Sample : K1817-14
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PST-027-B116-030619

Manual Integrations
 APPROVED

Sohil
 3/13/2019 4:33:43 PM

Quant Time: Mar 13 12:32:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 13 03:42:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	188308	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	642653	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	289147	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	520644	20.00	ng	0.00
76) Chrysene-d12	13.86	240	536922	20.00	ng	0.00
87) Perylene-d12	15.27	264	422764	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	876201	72.38	ng	0.00
7) Phenol-d6	6.37	99	1127911	74.17	ng	0.00
23) Nitrobenzene-d5	7.29	82	643436	59.91	ng	0.00
42) 2,4,6-Tribromophenol	10.55	330	299138	97.45	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	1220314	65.90	ng	0.00
79) Terphenyl-d14	12.82	244	1198879	43.28	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.48	163	107548	5.044	ng	98
71) Phenanthrene	11.26	178	86461	3.338	ng	100
75) Fluoranthene	12.44	202	143980	5.188	ng	97
78) Pyrene	12.67	202	181338	4.006	ng	98
83) Chrysene	13.89	228	87785	2.408	ng	97
88) Benzo(b)fluoranthene	14.87	252	69899m	2.556	ng	
90) Benzo(a)pyrene	15.21	252	51699	2.137	ng	# 93

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

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