

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031519\
 Data File : BF113065.D
 Acq On : 15 Mar 2019 14:58
 Operator : JU/SJ
 Sample : K1911-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RBR200036

Manual Integrations
 APPROVED

Sohil
 3/18/2019 2:59:29 PM

Quant Time: Mar 15 23:21:09 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	127271	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	513983	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	251808	20.00	ng	-0.01
64) Phenanthrene-d10	11.23	188	512801	20.00	ng	-0.01
76) Chrysene-d12	13.86	240	442883	20.00	ng	0.00
87) Perylene-d12	15.27	264	434875	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	886991	108.41	ng	0.00
7) Phenol-d6	6.37	99	1163217	113.18	ng	0.00
23) Nitrobenzene-d5	7.29	82	738312	85.95	ng	-0.01
42) 2,4,6-Tribromophenol	10.55	330	405612	151.73	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	1381395	93.29	ng	0.00
79) Terphenyl-d14	12.82	244	1544848	67.62	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.47	163	108356	5.836	ng	98
71) Phenanthrene	11.26	178	100551	3.941	ng	98
75) Fluoranthene	12.44	202	167069	6.112	ng	96
78) Pyrene	12.67	202	156179	4.183	ng	97
83) Chrysene	13.89	228	63325	2.106	ng	96
88) Benzo(b)fluoranthene	14.87	252	70682m	2.513	ng	

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

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