

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031519\
 Data File : BF113063.D
 Acq On : 15 Mar 2019 14:03
 Operator : JU/SJ
 Sample : K1922-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ORG-031419

Manual Integrations
 APPROVED

Sohil
 3/18/2019 2:59:28 PM

Quant Time: Mar 15 23:10:43 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	170089	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	632750	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	318455	20.00	ng	-0.01
64) Phenanthrene-d10	11.23	188	667548	20.00	ng	-0.01
76) Chrysene-d12	13.86	240	524340	20.00	ng	0.00
87) Perylene-d12	15.27	264	521371	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1171712	107.16	ng	0.00
7) Phenol-d6	6.37	99	1524783	111.01	ng	0.00
23) Nitrobenzene-d5	7.29	82	921006	87.10	ng	-0.01
42) 2,4,6-Tribromophenol	10.55	330	493832	146.07	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	1558557	79.78	ng	0.00
79) Terphenyl-d14	12.82	244	1798893	66.51	ng	0.00
Target Compounds						
10) Phenol	6.39	94	49833	3.109	ng	91
50) Dimethylphthalate	9.47	163	121252	5.164	ng	99
75) Fluoranthene	12.44	202	124164	3.489	ng	97
78) Pyrene	12.67	202	110555	2.501	ng	97
88) Benzo(b)fluoranthene	14.87	252	92037m	2.729	ng	
90) Benzo(a)pyrene	15.21	252	65945	2.211	ng	99

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

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