

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF031919\
 Data File : BF113149.D
 Acq On : 20 Mar 2019 3:06
 Operator : JU/SJ
 Sample : K1933-02
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-2

Manual Integrations
 APPROVED

Sohil
 3/20/2019 6:51:30 PM

Quant Time: Mar 20 07:29:22 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	6.72	152	191889	20.00	ng	-0.01
21) Naphthalene-d8	8.00	136	639521	20.00	ng	-0.01
39) Acenaphthene-d10	9.76	164	345398	20.00	ng	-0.01
64) Phenanthrene-d10	11.23	188	552565	20.00	ng	-0.01
76) Chrysene-d12	13.86	240	425171	20.00	ng	0.00
87) Perylene-d12	15.27	264	321637	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1133891	91.92	ng	0.00
7) Phenol-d6	6.37	99	1411362	91.08	ng	0.00
23) Nitrobenzene-d5	7.29	82	790387	73.95	ng	-0.01
42) 2,4,6-Tribromophenol	10.55	330	346574	94.52	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	1317621	58.15	ng	-0.01
79) Terphenyl-d14	12.82	244	1155130	52.67	ng	0.00
Target Compounds						
10) Phenol	6.39	94	67274	3.720	ng	91
50) Dimethylphthalate	9.47	163	242871	9.536	ng	99
71) Phenanthrene	11.26	178	77202	2.808	ng	98
75) Fluoranthene	12.44	202	176219	5.983	ng	96
78) Pyrene	12.67	202	148580	4.145	ng	97
81) Benzo(a)anthracene	13.85	228	87225	2.960	ng	99
83) Chrysene	13.89	228	66317	2.298	ng	97
88) Benzo(b)fluoranthene	14.87	252	85522m	4.111	ng	
90) Benzo(a)pyrene	15.21	252	53243	2.893	ng	99

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

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