

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031319\
 Data File : BF113043.D
 Acq On : 13 Mar 2019 22:13
 Operator : JU/SJ
 Sample : K1870-09 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-3

Manual Integrations
 APPROVED

Sohil
 3/14/2019 10:50:12 AM

Quant Time: Mar 14 02:38:39 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	198482	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	752564	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	368674	20.00	ng	-0.01
64) Phenanthrene-d10	11.23	188	680965	20.00	ng	-0.01
76) Chrysene-d12	13.86	240	566538	20.00	ng	0.00
87) Perylene-d12	15.27	264	526134	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	589053	46.17	ng	0.00
7) Phenol-d6	6.37	99	829022	51.72	ng	0.00
23) Nitrobenzene-d5	7.28	82	466803	37.12	ng	-0.02
42) 2,4,6-Tribromophenol	10.54	330	247493	63.23	ng	-0.01
45) 2-Fluorobiphenyl	9.08	172	813051	30.58	ng	-0.01
79) Terphenyl-d14	12.82	244	776497	26.57	ng	0.00
Target Compounds						
10) Phenol	6.38	94	55187	2.951	ng	92
50) Dimethylphthalate	9.47	163	87819	3.231	ng	100
71) Phenanthrene	11.26	178	127973	3.777	ng	96
75) Fluoranthene	12.44	202	311719	8.587	ng	96
78) Pyrene	12.67	202	310391	6.499	ng	99
81) Benzo(a)anthracene	13.85	228	154190	3.927	ng	98
83) Chrysene	13.89	228	147376	3.832	ng	97
86) Indeno(1,2,3-cd)pyrene	16.62	276	67923	2.071	ng	94
88) Benzo(b)fluoranthene	14.87	252	209993m	6.170	ng	
90) Benzo(a)pyrene	15.20	252	140153	4.656	ng	97
92) Benzo(a,h,i)perylene	17.03	276	68955	2.673	ng	98

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

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