

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030619\
 Data File : BF112900.D
 Acq On : 7 Mar 2019 00:15
 Operator : JU/SJ
 Sample : K1705-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EXT-028-SW004-022719

Manual Integrations
 APPROVED

Sohil
 3/7/2019 11:02:40 AM

Quant Time: Mar 07 05:29:18 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 01 15:40:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	204494	20.00	ng	-0.01
21) Naphthalene-d8	8.01	136	731622	20.00	ng	-0.01
39) Acenaphthene-d10	9.76	164	303641	20.00	ng	-0.01
64) Phenanthrene-d10	11.23	188	555205	20.00	ng	-0.01
76) Chrysene-d12	13.86	240	546723	20.00	ng	-0.01
87) Perylene-d12	15.26	264	592301	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	1147993	87.33	ng	0.00
7) Phenol-d6	6.37	99	1430580	86.63	ng	0.00
23) Nitrobenzene-d5	7.29	82	861938	70.50	ng	-0.02
42) 2,4,6-Tribromophenol	10.55	330	343237	106.48	ng	-0.01
45) 2-Fluorobiphenyl	9.09	172	1360318	71.09	ng	-0.01
79) Terphenyl-d14	12.82	244	1306289	46.32	ng	-0.01
Target Compounds						
32) Benzoic acid	7.72	122	15208	2.059	ng	88
50) Dimethylphthalate	9.48	163	95912	4.284	ng	98
71) Phenanthrene	11.26	178	62876	2.276	ng	98
75) Fluoranthene	12.44	202	203233	6.867	ng	98
78) Pyrene	12.67	202	180085	3.907	ng	98
81) Benzo(a)anthracene	13.85	228	95323	2.516	ng	95
83) Chrysene	13.89	228	134884	3.634	ng	96
84) Bis(2-ethylhexyl)phthalate	13.86	149	486693	20.395	ng	99
86) Indeno(1,2,3-cd)pyrene	16.62	276	63621m	2.011	ng	
88) Benzo(b)fluoranthene	14.87	252	182186m	4.755	ng	
89) Benzo(k)fluoranthene	14.89	252	69916m	2.015	ng	
90) Benzo(a)pyrene	15.20	252	107648	3.177	ng	98
92) Benzo(a,h,i)perylene	17.02	276	59209	2.039	ng	97

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

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