

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011819\
 Data File : BF112130.D
 Acq On : 18 Jan 2019 17:54
 Operator : JU/SJ
 Sample : K1191-04 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SITE4-EAST

Manual Integrations
 APPROVED

Sohil
 1/21/2019 12:00:50 PM

Quant Time: Jan 19 00:09:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 18 23:33:58 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	263275	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	1061369	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	537159	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	976790	20.00	ng	0.00
76) Chrysene-d12	14.02	240	625294	20.00	ng	0.00
87) Perylene-d12	15.50	264	602522	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	247774	17.08	ng	0.00
7) Phenol-d6	6.50	99	911762	50.24	ng	0.00
23) Nitrobenzene-d5	7.41	82	579349	37.71	ng	-0.02
42) 2,4,6-Tribromophenol	10.68	330	8932	1.54	ng	-0.01
45) 2-Fluorobiphenyl	9.21	172	1287948	35.15	ng	0.00
79) Terphenyl-d14	12.96	244	1162158	39.54	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.60	163	101950	2.602	ng	Qvalue # 98
71) Phenanthrene	11.40	178	422437	8.593	ng	97
72) Anthracene	11.45	178	102204	2.052	ng	97
75) Fluoranthene	12.59	202	440893	8.495	ng	99
78) Pyrene	12.82	202	371608	9.008	ng	98
81) Benzo(a)anthracene	14.02	228	149567	4.172	ng	99
83) Chrysene	14.04	228	136697	4.029	ng	96
88) Benzo(b)fluoranthene	15.07	252	149672m	4.376	ng	
90) Benzo(a)pyrene	15.44	252	108624	3.463	ng	# 55
92) Benzo(g,h,i)perylene	17.43	276	55807	2.033	ng	# 79

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

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