

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011119\
 Data File : BF112020.D
 Acq On : 11 Jan 2019 17:28
 Operator : JU/SJ
 Sample : K1062-01RE
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-(1-10)-COMP

Manual Integrations
 APPROVED

Sohil
 1/14/2019 4:10:18 PM

Quant Time: Jan 11 22:58:32 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 11 21:37:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	104981	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	378345	20.00	ng	0.00
39) Acenaphthene-d10	9.91	164	169669	20.00	ng	0.00
64) Phenanthrene-d10	11.40	188	283032	20.00	ng	0.00
76) Chrysene-d12	14.04	240	270286	20.00	ng	0.00
87) Perylene-d12	15.53	264	191746	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	513160	83.63	ng	0.00
7) Phenol-d6	6.52	99	659252	83.35	ng	0.00
23) Nitrobenzene-d5	7.43	82	421182	59.50	ng	-0.01
42) 2,4,6-Tribromophenol	10.70	330	160742	72.84	ng	0.00
45) 2-Fluorobiphenyl	9.23	172	665749	57.22	ng	0.00
79) Terphenyl-d14	12.98	244	594540	41.75	ng	0.00
Target Compounds						
10) Phenol	6.53	94	29716	3.457	ng	Qvalue # 76
50) Dimethylphthalate	9.62	163	68810	5.399	ng	99
71) Phenanthrene	11.42	178	262368	17.338	ng	98
72) Anthracene	11.47	178	49096	3.196	ng	97
73) Carbazole	11.62	167	31988	2.141	ng	99
75) Fluoranthene	12.61	202	381144	24.535	ng	99
78) Pyrene	12.84	202	319641	17.187	ng	99
80) Butylbenzylphthalate	13.45	149	24154	3.072	ng	94
81) Benzo(a)anthracene	14.03	228	186015	11.958	ng	99
83) Chrysene	14.06	228	181506	12.238	ng	96
84) Bis(2-ethylhexyl)phthalate	14.01	149	54124	5.242	ng	99
86) Indeno(1,2,3-cd)pyrene	17.01	276	59906	5.540	ng	96
88) Benzo(b)fluoranthene	15.09	252	184236m	16.062	ng	
89) Benzo(k)fluoranthene	15.12	252	57899m	5.370	ng	
90) Benzo(a)pyrene	15.46	252	105689	10.479	ng	98
91) Dibenzo(a,h)anthracene	17.02	278	16894	2.083	ng	# 78
92) Benzo(a,h,i)perylene	17.46	276	58740	7.404	ng	# 94

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

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