

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF103020\
 Data File : BF122197.D
 Acq On : 30 Oct 2020 23:56
 Operator : JU/CG
 Sample : L4564-03
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-03-017-B

Manual Integrations
 APPROVED

mohammad
 11/2/2020 12:36:26 PM

Quant Time: Nov 02 07:48:33 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 30 10:16:54 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.710	152	96860	20.00	ng	0.00
21) Naphthalene-d8	7.992	136	371032	20.00	ng	0.00
39) Acenaphthene-d10	9.745	164	178179	20.00	ng	0.00
64) Phenanthrene-d10	11.233	188	274718	20.00	ng	0.00
76) Chrysene-d12	13.880	240	190543	20.00	ng	0.00
86) Perylene-d12	15.315	264	197708	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.345	112	238326	36.32	ng	0.01
7) Phenol-d6	6.369	99	421184	49.86	ng	0.00
23) Nitrobenzene-d5	7.281	82	329858	45.59	ng	0.00
42) 2,4,6-Tribromophenol	10.545	330	45248	20.53	ng	0.00
45) 2-Fluorobiphenyl	9.069	172	577975	47.73	ng	0.00
79) Terphenyl-d14	12.816	244	394835	39.49	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.387	94	26906	3.09	ng	# 81
50) Dimethylphthalate	9.469	163	112564	8.64	ng	100
71) Phenanthrene	11.257	178	41081	2.73	ng	99
75) Fluoranthene	12.451	202	85767	5.31	ng	98
78) Pyrene	12.680	202	79572	5.48	ng	99
81) Benzo(a)anthracene	13.868	228	44495	3.56	ng	98
83) Chrysene	13.910	228	42573	3.47	ng	98
87) Indeno(1,2,3-cd)pyrene	16.751	276	31198m	2.43	ng	
88) Benzo(b)fluoranthene	14.910	252	65047m	4.87	ng	
90) Benzo(a)pyrene	15.262	252	46719	4.05	ng	# 96
92) Benzo(g,h,i)perylene	17.180	276	30190	2.85	ng	97

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

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