

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071020\
 Data File : BF120848.D
 Acq On : 10 Jul 2020 23:31
 Operator : JU/CG
 Sample : L3184-01 5X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-02-070920

Manual Integrations
 APPROVED

mohammad
 7/13/2020 1:05:14 PM

Quant Time: Jul 11 03:28:42 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 02 18:51:52 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	180874	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	651444	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	304241	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	471791	20.00	ng	0.00
76) Chrysene-d12	14.01	240	379234	20.00	ng	0.00
86) Perylene-d12	15.47	264	447542	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	243970	20.82	ng	0.00
7) Phenol-d6	6.46	99	287962	19.24	ng	-0.01
23) Nitrobenzene-d5	7.40	82	170181	15.14	ng	-0.01
42) 2,4,6-Tribromophenol	10.67	330	49430	16.17	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	342216	17.02	ng	0.00
79) Terphenyl-d14	12.96	244	276074	14.25	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.60	163	58006	2.614	ng	Qvalue # 98
71) Phenanthrene	11.39	178	106308	4.153	ng	99
75) Fluoranthene	12.58	202	202879	7.440	ng	99
78) Pyrene	12.81	202	171053	5.648	ng	98
81) Benzo(a)anthracene	14.00	228	98856	4.067	ng	97
83) Chrysene	14.03	228	77212	3.114	ng	96
88) Benzo(b)fluoranthene	15.04	252	112565m	3.826	ng	
90) Benzo(a)pyrene	15.41	252	84525	3.220	ng	# 94
92) Benzo(g,h,i)perylene	17.36	276	47060	2.152	ng	# 91

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

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