

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071420\
 Data File : BF120866.D
 Acq On : 14 Jul 2020 17:49
 Operator : JU/CG
 Sample : L3287-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-03-012-A

Manual Integrations
 APPROVED

mohammad
 7/15/2020 10:59:52 AM

Quant Time: Jul 14 18:19:38 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 13 14:16:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	196754	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	722707	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	335301	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	537385	20.00	ng	0.00
76) Chrysene-d12	13.99	240	458858	20.00	ng	-0.01
86) Perylene-d12	15.46	264	448592	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	749147	58.77	ng	0.00
7) Phenol-d6	6.46	99	896493	55.08	ng	0.00
23) Nitrobenzene-d5	7.39	82	546616	43.82	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	193599	57.47	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	999046	45.08	ng	0.00
79) Terphenyl-d14	12.94	244	802304	34.23	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.58	163	163443	6.682	ng	98
71) Phenanthrene	11.37	178	94182	3.230	ng	99
75) Fluoranthene	12.56	202	223824	7.206	ng	98
78) Pyrene	12.79	202	211106	5.761	ng	98
81) Benzo(a)anthracene	13.98	228	141938m	4.826	ng	
83) Chrysene	14.02	228	131429	4.381	ng	96
87) Indeno(1,2,3-cd)pyrene	16.89	276	76421	2.656	ng	95
88) Benzo(b)fluoranthene	15.03	252	189274m	6.418	ng	
89) Benzo(k)fluoranthene	15.05	252	64327m	2.242	ng	
90) Benzo(a)pyrene	15.39	252	118892	4.519	ng	# 94
92) Benzo(g,h,i)perylene	17.33	276	72566	3.311	ng	97

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

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