

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082120\
 Data File : BF121379.D
 Acq On : 21 Aug 2020 21:20
 Operator : JU/CG
 Sample : L3734-06 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EROC-BF-1E

Manual Integrations
 APPROVED

Sohil
 8/24/2020 1:50:09 PM

Quant Time: Aug 31 09:02:50 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF081920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 20 04:56:10 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	164210	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	630335	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	307885	20.00	ng	0.00
64) Phenanthrene-d10	11.21	188	527813	20.00	ng	0.00
76) Chrysene-d12	13.85	240	340649	20.00	ng	0.00
86) Perylene-d12	15.26	264	392775	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	84966	8.48	ng	0.00
7) Phenol-d6	6.33	99	319390	24.98	ng	0.00
23) Nitrobenzene-d5	7.26	82	316393	28.39	ng	0.00
42) 2,4,6-Tribromophenol	10.52	330	10033	2.80	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	611621	29.73	ng	0.00
79) Terphenyl-d14	12.79	244	480795	27.58	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.45	163	55224	2.465	ng	99
71) Phenanthrene	11.23	178	132817	4.862	ng	99
75) Fluoranthene	12.42	202	188733	6.365	ng	99
78) Pyrene	12.65	202	162060	6.333	ng	99
81) Benzo(a)anthracene	13.83	228	78603	3.715	ng	# 93
83) Chrysene	13.87	228	72602	3.355	ng	# 95
88) Benzo(b)fluoranthene	14.86	252	92028m	3.627	ng	
90) Benzo(a)pyrene	15.20	252	69235	3.158	ng	# 88
92) Benzo(g,h,i)perylene	17.03	276	47299	2.461	ng	# 91

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

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