

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082120\
 Data File : BF121381.D
 Acq On : 21 Aug 2020 22:22
 Operator : JU/CG
 Sample : L3734-04 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EROC-BF-1C

Manual Integrations
 APPROVED

Sohil
 8/24/2020 1:50:11 PM

Quant Time: Aug 31 09:02:08 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	163613	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	621020	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	300362	20.00	ng	0.00
64) Phenanthrene-d10	11.21	188	500827	20.00	ng	0.00
76) Chrysene-d12	13.85	240	336914	20.00	ng	0.00
86) Perylene-d12	15.27	264	377437	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	104421	10.47	ng	0.00
7) Phenol-d6	6.33	99	446838	35.08	ng	0.00
23) Nitrobenzene-d5	7.26	82	323296	29.44	ng	0.00
42) 2,4,6-Tribromophenol	10.52	330	7078	2.02	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	605626	30.40	ng	0.00
79) Terphenyl-d14	12.80	244	450626	26.14	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.45	163	59775	2.735	ng	Qvalue 100
71) Phenanthrene	11.23	178	169896	6.555	ng	99
75) Fluoranthene	12.42	202	224159	7.967	ng	98
78) Pyrene	12.65	202	211899	8.373	ng	100
81) Benzo(a)anthracene	13.84	228	109621	5.238	ng	95
83) Chrysene	13.87	228	99387	4.644	ng	97
84) Bis(2-ethylhexyl)phthalate	13.83	149	82954	6.487	ng	99
88) Benzo(b)fluoranthene	14.87	252	126629m	5.193	ng	
89) Benzo(k)fluoranthene	14.89	252	43580m	2.030	ng	
90) Benzo(a)pyrene	15.21	252	92991	4.414	ng	# 91
92) Benzo(g,h,i)perylene	17.04	276	54484	2.950	ng	# 89

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

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