

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082120\
 Data File : BF121383.D
 Acq On : 21 Aug 2020 23:23
 Operator : JU/CG
 Sample : L3734-02 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EROC-BF-1A

Manual Integrations
 APPROVED

Sohil
 8/24/2020 1:50:13 PM

Quant Time: Aug 22 03:00:09 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	6.70	152	159424	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	592758	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	271708	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	436879	20.00	ng	0.00
76) Chrysene-d12	13.85	240	337500	20.00	ng	0.00
86) Perylene-d12	15.27	264	349276	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	178448	18.35	ng	0.00
7) Phenol-d6	6.34	99	479472	38.63	ng	0.00
23) Nitrobenzene-d5	7.26	82	335839	32.04	ng	0.00
42) 2,4,6-Tribromophenol	10.52	330	21685	6.85	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	618843	36.32	ng	0.00
79) Terphenyl-d14	12.80	244	456123	26.41	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.45	163	57299	2.899	ng	98
71) Phenanthrene	11.23	178	86383	3.820	ng	99
75) Fluoranthene	12.42	202	146089	5.953	ng	97
78) Pyrene	12.65	202	130391	5.143	ng	99
81) Benzo(a)anthracene	13.84	228	68287	3.258	ng	# 92
83) Chrysene	13.88	228	72889	3.400	ng	# 95
84) Bis(2-ethylhexyl)phthalate	13.84	149	101603	7.931	ng	# 96
88) Benzo(b)fluoranthene	14.87	252	88325m	3.914	ng	
90) Benzo(a)pyrene	15.21	252	56297	2.888	ng	# 78

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

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