

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072420\
 Data File : BF121018.D
 Acq On : 24 Jul 2020 22:52
 Operator : JU/CG
 Sample : L3383-13
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CRT-022

Manual Integrations
 APPROVED

mohammad
 7/27/2020 12:19:54 PM

Quant Time: Jul 27 02:42:43 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 16 14:20:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	152379	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	541109	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	276597	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	465902	20.00	ng	0.00
76) Chrysene-d12	13.97	240	380472	20.00	ng	0.00
86) Perylene-d12	15.42	264	419421	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	755833	81.32	ng	0.00
7) Phenol-d6	6.45	99	890428	74.16	ng	0.00
23) Nitrobenzene-d5	7.37	82	557340	59.60	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	214507	70.85	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	1111284	59.99	ng	0.00
79) Terphenyl-d14	12.92	244	927842	53.01	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.56	163	97642	4.893	ng	99
71) Phenanthrene	11.36	178	184367	7.694	ng	98
75) Fluoranthene	12.55	202	269852	10.160	ng	96
78) Pyrene	12.77	202	250027	9.295	ng	99
81) Benzo(a)anthracene	13.96	228	145933m	6.334	ng	
83) Chrysene	13.99	228	139413	5.758	ng	97
84) Bis(2-ethylhexyl)phthalate	13.95	149	36164	2.446	ng	99
87) Indeno(1,2,3-cd)pyrene	16.84	276	80038	2.744	ng	96
88) Benzo(b)fluoranthene	15.00	252	209566m	8.266	ng	
89) Benzo(k)fluoranthene	15.02	252	64538m	2.791	ng	
90) Benzo(a)pyrene	15.35	252	136369	5.896	ng	96
92) Benzo(a,h,i)perylene	17.27	276	76256	3.246	ng	97

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

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