

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060220\
 Data File : BF120488.D
 Acq On : 2 Jun 2020 12:13
 Operator : JU/CG
 Sample : L2798-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM88-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/3/2020 4:35:47 PM

Quant Time: Jun 02 13:39:44 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 02 11:37:51 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	218755	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	803443	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	385303	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	644559	20.00	ng	0.00
76) Chrysene-d12	14.01	240	415130	20.00	ng	0.00
86) Perylene-d12	15.47	264	458049	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	877801	76.47	ng	0.00
7) Phenol-d6	6.48	99	1134343	80.18	ng	0.00
23) Nitrobenzene-d5	7.40	82	714345	58.16	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	287530	80.54	ng	-0.01
45) 2-Fluorobiphenyl	9.20	172	1343094	59.14	ng	0.00
79) Terphenyl-d14	12.95	244	1133509	60.76	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	235033	9.757	ng	99
71) Phenanthrene	11.39	178	324781	11.435	ng	98
72) Anthracene	11.44	178	57984	2.035	ng	98
75) Fluoranthene	12.58	202	377956	12.368	ng	97
78) Pyrene	12.81	202	354345	12.274	ng	98
81) Benzo(a)anthracene	14.00	228	165584	7.442	ng	98
83) Chrysene	14.03	228	164655	7.059	ng	98
87) Indeno(1,2,3-cd)pyrene	16.92	276	104476	4.170	ng	94
88) Benzo(b)fluoranthene	15.04	252	234247m	9.207	ng	
89) Benzo(k)fluoranthene	15.07	252	59290m	2.496	ng	
90) Benzo(a)pyrene	15.40	252	163366	7.341	ng	98
92) Benzo(a,h,i)perylene	17.36	276	96749	4.803	ng	100

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

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