

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060220\
 Data File : BF120489.D
 Acq On : 2 Jun 2020 12:40
 Operator : JU/CG
 Sample : L2798-09
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM99-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/3/2020 4:35:50 PM

Quant Time: Jun 02 13:46:32 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	242360	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	871022	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	414927	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	667851	20.00	ng	0.00
76) Chrysene-d12	14.01	240	440047	20.00	ng	0.00
86) Perylene-d12	15.47	264	478747	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	1112776	87.50	ng	0.01
7) Phenol-d6	6.48	99	1432756	91.40	ng	0.00
23) Nitrobenzene-d5	7.40	82	976412	73.33	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	351673	91.47	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	1694590	69.29	ng	0.00
79) Terphenyl-d14	12.96	244	1404952	71.05	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	307958	11.872	ng	99
71) Phenanthrene	11.39	178	222703	7.568	ng	99
72) Anthracene	11.44	178	59094	2.001	ng	97
75) Fluoranthene	12.58	202	443434	14.004	ng	98
78) Pyrene	12.81	202	395394	12.920	ng	98
81) Benzo(a)anthracene	14.00	228	211188	8.955	ng	97
83) Chrysene	14.03	228	204679m	8.278	ng	
87) Indeno(1,2,3-cd)pyrene	16.92	276	121967	4.658	ng	96
88) Benzo(b)fluoranthene	15.04	252	304614m	11.456	ng	
89) Benzo(k)fluoranthene	15.07	252	80957m	3.261	ng	
90) Benzo(a)pyrene	15.40	252	203222	8.737	ng	99
92) Benzo(a,h,i)perylene	17.36	276	112640	5.350	ng	99

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

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