

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060320\
 Data File : BF120518.D
 Acq On : 3 Jun 2020 19:15
 Operator : JU/CG
 Sample : L2839-20
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM21-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/4/2020 12:12:43 PM

Quant Time: Jun 04 01:12:52 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 03 14:26:17 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	144353	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	533061	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	264653	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	437523	20.00	ng	0.00
76) Chrysene-d12	14.00	240	421331	20.00	ng	-0.01
86) Perylene-d12	15.45	264	386168	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	596547	78.76	ng	0.00
7) Phenol-d6	6.47	99	765793	82.02	ng	0.00
23) Nitrobenzene-d5	7.40	82	532012	65.28	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	222521	90.75	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	1034014	66.29	ng	0.00
79) Terphenyl-d14	12.94	244	1117844	59.04	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	157220	9.502	ng	98
71) Phenanthrene	11.38	178	184641	9.578	ng	97
75) Fluoranthene	12.57	202	323370	15.589	ng	99
78) Pyrene	12.80	202	298025	10.171	ng	97
81) Benzo(a)anthracene	13.99	228	166882	7.390	ng	99
83) Chrysene	14.02	228	140081	5.917	ng	96
87) Indeno(1,2,3-cd)pyrene	16.89	276	58191	2.755	ng	95
88) Benzo(b)fluoranthene	15.03	252	173708m	8.099	ng	
89) Benzo(k)fluoranthene	15.05	252	60463m	3.020	ng	
90) Benzo(a)pyrene	15.39	252	118314	6.306	ng	99
92) Benzo(g,h,i)perylene	17.33	276	60675	3.573	ng	96

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

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