

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060420\
 Data File : BF120549.D
 Acq On : 4 Jun 2020 22:33
 Operator : JU/CG
 Sample : L2839-07RE 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM7-COMP-2020-0-6RE

Manual Integrations
 APPROVED

mohammad
 6/5/2020 11:58:16 AM

Quant Time: Jun 05 04:33:35 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 04 16:38:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	210074	20.00	ng	0.00
21) Naphthalene-d8	8.11	136	769315	20.00	ng	0.00
39) Acenaphthene-d10	9.86	164	275994	20.00	ng	0.00
64) Phenanthrene-d10	11.35	188	479193	20.00	ng	0.00
76) Chrysene-d12	14.00	240	440021	20.00	ng	0.00
86) Perylene-d12	15.46	264	310769	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	493522	44.77	ng	0.00
7) Phenol-d6	6.46	99	619200	45.57	ng	-0.01
23) Nitrobenzene-d5	7.39	82	426754	36.29	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	138866	54.30	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	643874	39.58	ng	0.00
79) Terphenyl-d14	12.94	244	751360	38.00	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.58	163	97981	5.679	ng	99
71) Phenanthrene	11.37	178	120387	5.702	ng	97
75) Fluoranthene	12.57	202	245404	10.801	ng	97
78) Pyrene	12.80	202	284562	9.299	ng	98
81) Benzo(a)anthracene	13.99	228	146298	6.204	ng	97
83) Chrysene	14.02	228	138729	5.611	ng	96
87) Indeno(1,2,3-cd)pyrene	16.90	276	56040	3.297	ng	97
88) Benzo(b)fluoranthene	15.03	252	196045m	11.358	ng	
89) Benzo(k)fluoranthene	15.06	252	43411m	2.694	ng	
90) Benzo(a)pyrene	15.39	252	90607	6.001	ng	100
92) Benzo(g,h,i)perylene	17.34	276	52861	3.868	ng	97

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

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