

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060320\
 Data File : BF120507.D
 Acq On : 3 Jun 2020 14:13
 Operator : JU/CG
 Sample : L2839-21
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM34-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/4/2020 12:12:11 PM

Quant Time: Jun 04 00:52:36 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	234097	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	887408	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	455941	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	841437	20.00	ng	0.00
76) Chrysene-d12	14.00	240	540534	20.00	ng	-0.01
86) Perylene-d12	15.45	264	535864	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	797844	64.95	ng	0.00
7) Phenol-d6	6.46	99	1045735	69.07	ng	-0.01
23) Nitrobenzene-d5	7.39	82	772163	56.92	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	339644	80.40	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	1446816	53.84	ng	0.00
79) Terphenyl-d14	12.94	244	1482378	61.03	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.58	163	213426	7.488	ng	100
71) Phenanthrene	11.38	178	156817	4.230	ng	96
75) Fluoranthene	12.57	202	314037	7.872	ng	98
78) Pyrene	12.80	202	281485	7.488	ng	97
81) Benzo(a)anthracene	13.99	228	121419	4.191	ng	96
83) Chrysene	14.02	228	120656	3.973	ng	97
87) Indeno(1,2,3-cd)pyrene	16.89	276	78221	2.669	ng	94
88) Benzo(b)fluoranthene	15.03	252	158566m	5.328	ng	
90) Benzo(a)pyrene	15.39	252	108050	4.150	ng	98
92) Benzo(g,h,i)perylene	17.33	276	74182	3.148	ng	98

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

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