

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060320\
 Data File : BF120527.D
 Acq On : 3 Jun 2020 23:22
 Operator : JU/CG
 Sample : L2839-07 2X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

mohammad
 6/4/2020 12:13:09 PM

Quant Time: Jun 04 02:24:13 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	134042	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	499982	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	293221	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	486890	20.00	ng	0.00
76) Chrysene-d12	14.00	240	394206	20.00	ng	-0.01
86) Perylene-d12	15.46	264	325586	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	447106	63.57	ng	0.00
7) Phenol-d6	6.47	99	457491	52.77	ng	0.00
23) Nitrobenzene-d5	7.39	82	265399	34.72	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	138444	50.96	ng	-0.01
45) 2-Fluorobiphenyl	9.19	172	703328	40.70	ng	0.00
79) Terphenyl-d14	12.94	244	478331	27.00	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.59	163	106134	5.790	ng	98
71) Phenanthrene	11.38	178	117474	5.476	ng	97
75) Fluoranthene	12.57	202	250864	10.867	ng	98
78) Pyrene	12.80	202	232660	8.487	ng	98
81) Benzo(a)anthracene	13.99	228	139986	6.626	ng	98
83) Chrysene	14.02	228	115776	5.227	ng	97
87) Indeno(1,2,3-cd)pyrene	16.90	276	60007	3.370	ng	99
88) Benzo(b)fluoranthene	15.03	252	147021m	8.130	ng	
89) Benzo(k)fluoranthene	15.06	252	37543m	2.224	ng	
90) Benzo(a)pyrene	15.39	252	96970	6.130	ng	96
92) Benzo(g,h,i)perylene	17.34	276	59632	4.164	ng	95

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

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