

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060320\
 Data File : BF120512.D
 Acq On : 3 Jun 2020 16:30
 Operator : JU/CG
 Sample : L2839-15
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM104-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/4/2020 12:12:26 PM

Quant Time: Jun 03 17:17:59 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	221738	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	833733	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	409900	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	692162	20.00	ng	0.00
76) Chrysene-d12	14.00	240	464008	20.00	ng	-0.01
86) Perylene-d12	15.45	264	381289	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	840279	72.22	ng	0.00
7) Phenol-d6	6.47	99	1227961	85.62	ng	0.00
23) Nitrobenzene-d5	7.40	82	803675	63.05	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	335941	88.45	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	1359307	56.27	ng	0.00
79) Terphenyl-d14	12.94	244	1302866	62.48	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	317209	12.378	ng	99
71) Phenanthrene	11.38	178	80975	2.655	ng	97
75) Fluoranthene	12.57	202	175427	5.346	ng	98
78) Pyrene	12.80	202	167121	5.179	ng	97
81) Benzo(a)anthracene	13.99	228	88288	3.550	ng	96
83) Chrysene	14.02	228	86354m	3.312	ng	
87) Indeno(1,2,3-cd)pyrene	16.89	276	50236	2.409	ng	96
88) Benzo(b)fluoranthene	15.03	252	94805m	4.477	ng	
90) Benzo(a)pyrene	15.39	252	66196	3.573	ng	97
92) Benzo(g,h,i)perylene	17.33	276	48454	2.889	ng	96

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

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