

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060420\
 Data File : BF120538.D
 Acq On : 4 Jun 2020 17:24
 Operator : JU/CG
 Sample : L2839-19
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM20-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/5/2020 11:57:37 AM

Quant Time: Jun 04 18:59:27 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	163869	20.00	ng	0.00
21) Naphthalene-d8	8.11	136	547159	20.00	ng	0.00
39) Acenaphthene-d10	9.86	164	370562	20.00	ng	0.00
64) Phenanthrene-d10	11.35	188	608391	20.00	ng	0.00
76) Chrysene-d12	13.99	240	469045	20.00	ng	-0.01
86) Perylene-d12	15.45	264	442411	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	857780	99.76	ng	0.00
7) Phenol-d6	6.46	99	1015863	95.85	ng	0.00
23) Nitrobenzene-d5	7.39	82	712347	85.16	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	330817	96.35	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	1429603	65.46	ng	0.00
79) Terphenyl-d14	12.95	244	1462987	69.41	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.58	163	291063	12.564	ng	Qvalue 100
71) Phenanthrene	11.37	178	114331	4.265	ng	98
75) Fluoranthene	12.56	202	277938	9.635	ng	99
78) Pyrene	12.79	202	263907	8.090	ng	98
81) Benzo(a)anthracene	13.98	228	143084	5.692	ng	97
83) Chrysene	14.02	228	137687	5.224	ng	97
87) Indeno(1,2,3-cd)pyrene	16.89	276	72282	2.987	ng	97
88) Benzo(b)fluoranthene	15.03	252	188145m	7.657	ng	
89) Benzo(k)fluoranthene	15.05	252	51413m	2.241	ng	
90) Benzo(a)pyrene	15.39	252	118303	5.504	ng	99
92) Benzo(g,h,i)perylene	17.33	276	67127	3.450	ng	98

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

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