

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060520\
 Data File : BF120555.D
 Acq On : 5 Jun 2020 13:10
 Operator : JU/CG
 Sample : L2839-17
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM108-2-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/8/2020 12:50:50 PM

Quant Time: Jun 06 02:58:22 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 05 15:36:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	203473	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	802833	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	393347	20.00	ng	0.00
64) Phenanthrene-d10	11.34	188	558429	20.00	ng	0.00
76) Chrysene-d12	13.97	240	361011	20.00	ng	0.00
86) Perylene-d12	15.43	264	362277	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	985275	92.28	ng	0.00
7) Phenol-d6	6.45	99	1341752	101.96	ng	0.00
23) Nitrobenzene-d5	7.38	82	862774	70.30	ng	0.00
42) 2,4,6-Tribromophenol	10.65	330	276554	75.88	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	1515757	65.38	ng	0.00
79) Terphenyl-d14	12.93	244	1193035	73.54	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.57	163	240310	9.772	ng	99
71) Phenanthrene	11.36	178	60603	2.463	ng	99
75) Fluoranthene	12.55	202	179419	6.777	ng	94
78) Pyrene	12.78	202	166804	6.644	ng	99
81) Benzo(a)anthracene	13.96	228	88181	4.558	ng	99
83) Chrysene	14.00	228	82037	4.044	ng	97
87) Indeno(1,2,3-cd)pyrene	16.85	276	68925	3.479	ng	99
88) Benzo(b)fluoranthene	15.00	252	108557m	5.395	ng	
90) Benzo(a)pyrene	15.36	252	75026	4.263	ng	98
92) Benzo(g,h,i)perylene	17.29	276	66736	4.189	ng	# 94

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

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