

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060420\
 Data File : BF120535.D
 Acq On : 4 Jun 2020 16:01
 Operator : JU/CG
 Sample : L2839-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM31-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/5/2020 11:57:26 AM

Quant Time: Jun 04 17:00:12 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	175980	20.00	ng	0.00
21) Naphthalene-d8	8.11	136	664005	20.00	ng	-0.01
39) Acenaphthene-d10	9.86	164	341381	20.00	ng	-0.02
64) Phenanthrene-d10	11.35	188	623791	20.00	ng	-0.01
76) Chrysene-d12	13.99	240	383379	20.00	ng	-0.02
86) Perylene-d12	15.45	264	428251	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	971471	105.20	ng	0.00
7) Phenol-d6	6.46	99	1225490	107.67	ng	-0.01
23) Nitrobenzene-d5	7.39	82	787700	77.60	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	348439	110.16	ng	-0.01
45) 2-Fluorobiphenyl	9.19	172	1439996	71.57	ng	-0.01
79) Terphenyl-d14	12.94	244	1322361	76.75	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.58	163	285776	13.390	ng	99
71) Phenanthrene	11.37	178	81703	2.973	ng	95
75) Fluoranthene	12.56	202	176442	5.966	ng	98
78) Pyrene	12.79	202	166024	6.227	ng	98
81) Benzo(a)anthracene	13.98	228	79969	3.892	ng	98
83) Chrysene	14.02	228	72569	3.369	ng	95
87) Indeno(1,2,3-cd)pyrene	16.89	276	51717	2.208	ng	95
88) Benzo(b)fluoranthene	15.02	252	112760m	4.741	ng	
90) Benzo(a)pyrene	15.39	252	79821	3.836	ng	97
92) Benzo(g,h,i)perylene	17.33	276	44249	2.349	ng	98

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

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