

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060420\
 Data File : BF120548.D
 Acq On : 4 Jun 2020 22:06
 Operator : JU/CG
 Sample : L2839-08RE 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM112-COMP-2020-0-6RE

Manual Integrations
 APPROVED

mohammad
 6/5/2020 11:58:14 AM

Quant Time: Jun 05 01:31:10 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	216530	20.00	ng	0.00
21) Naphthalene-d8	8.11	136	818590	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	273633	20.00	ng	0.00
64) Phenanthrene-d10	11.35	188	445451	20.00	ng	0.00
76) Chrysene-d12	14.00	240	427735	20.00	ng	0.00
86) Perylene-d12	15.46	264	339070	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	631364	55.57	ng	0.00
7) Phenol-d6	6.46	99	742604	53.03	ng	0.00
23) Nitrobenzene-d5	7.39	82	467702	37.37	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	141745	55.91	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	701728	43.51	ng	0.00
79) Terphenyl-d14	12.94	244	633220	32.94	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.58	163	114210	6.676	ng	99
71) Phenanthrene	11.37	178	121074	6.168	ng	98
75) Fluoranthene	12.57	202	249605	11.818	ng	96
78) Pyrene	12.80	202	232579	7.819	ng	97
81) Benzo(a)anthracene	13.99	228	160809	7.015	ng	96
83) Chrysene	14.02	228	152376m	6.340	ng	
87) Indeno(1,2,3-cd)pyrene	16.90	276	57342	3.092	ng	# 94
88) Benzo(b)fluoranthene	15.03	252	164358m	8.727	ng	
89) Benzo(k)fluoranthene	15.06	252	62733m	3.568	ng	
90) Benzo(a)pyrene	15.39	252	115382	7.004	ng	99
92) Benzo(g,h,i)perylene	17.34	276	54877	3.680	ng	98

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

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