

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100520\
 Data File : BF121836.D
 Acq On : 6 Oct 2020 1:00
 Operator : JU/CG
 Sample : L4247-02MS 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-18(0-1)MS

Manual Integrations
 APPROVED

mohammad
 10/7/2020 10:36:41 AM

Quant Time: Oct 06 02:54:27 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 01 14:29:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	87479	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	320629	20.00	ng	-0.01
39) Acenaphthene-d10	9.80	164	146062	20.00	ng	-0.01
64) Phenanthrene-d10	11.29	188	230112	20.00	ng	0.00
76) Chrysene-d12	13.92	240	235361	20.00	ng	-0.01
86) Perylene-d12	15.36	264	213168	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	14150	2.40	ng	0.00
7) Phenol-d6	6.41	99	18768	2.43	ng	0.00
23) Nitrobenzene-d5	7.33	82	14302	2.39	ng	-0.01
42) 2,4,6-Tribromophenol	10.59	330	3411	1.88	ng	-0.01
45) 2-Fluorobiphenyl	9.12	172	24900	2.58	ng	-0.01
79) Terphenyl-d14	12.86	244	27079	2.33	ng	-0.02
Target Compounds						
22) Acetophenone	7.16	105	18630	2.290	ng	# 46
26) 2-Nitrophenol	7.67	139	2190	3.241	ng	# 61
32) Benzoic acid	7.83	122	338	7.008	ng	# 64
35) Caprolactam	8.47	113	6496	3.991	ng	# 11
37) 2-Methylnaphthalene	8.76	142	28921	2.607	ng	98
38) 1-Methylnaphthalene	8.86	142	25002m	2.422	ng	
65) 4,6-Dinitro-2-methylphenol	10.61	198	581	7.205	ng	# 1
66) n-Nitrosodiphenylamine	10.46	169	16413	2.079	ng	# 85

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

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