

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081420\
 Data File : BG046291.D
 Acq On : 15 Aug 2020 1:25
 Operator : CG/JU
 Sample : L3660-04
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 7211979

Manual Integrations
 APPROVED

mohammad
 8/18/2020 1:41:51 PM

Quant Time: Aug 17 02:51:34 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 30 18:43:49 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.95	152	111695	20.00	ng	0.00
21) Naphthalene-d8	10.74	136	466119	20.00	ng	0.00
39) Acenaphthene-d10	14.57	164	302444	20.00	ng	0.00
64) Phenanthrene-d10	17.32	188	632051	20.00	ng	0.00
76) Chrysene-d12	21.57	240	600678	20.00	ng	0.00
86) Perylene-d12	24.68	264	644697	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	515537	80.24	ng	0.00
7) Phenol-d6	7.12	99	685697	78.30	ng	0.00
23) Nitrobenzene-d5	9.10	82	475969	55.29	ng	0.00
42) 2,4,6-Tribromophenol	16.07	330	365192	78.61	ng	0.00
45) 2-Fluorobiphenyl	13.20	172	1153508	55.40	ng	0.00
79) Terphenyl-d14	19.93	244	1371496	46.50	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.03	163	79573	3.428	ng	98
71) Phenanthrene	17.36	178	438489	12.665	ng	97
75) Fluoranthene	19.37	202	1177290	27.236	ng	99
78) Pyrene	19.73	202	1011022	25.050	ng	98
81) Benzo(a)anthracene	21.55	228	492411	12.250	ng	98
83) Chrysene	21.61	228	557660	14.162	ng	99
87) Indeno(1,2,3-cd)pyrene	28.19	276	409411	8.327	ng	96
88) Benzo(b)fluoranthene	23.70	252	805595	18.647	ng	99
89) Benzo(k)fluoranthene	23.76	252	266030m	6.278	ng	
90) Benzo(a)pyrene	24.53	252	553970	13.780	ng	99
91) Dibenzo(a,h)anthracene	28.24	278	85543	2.081	ng	95
92) Benzo(a,h,i)perylene	29.29	276	396226	9.545	ng	98

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

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