

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061620\
 Data File : BG045780.D
 Acq On : 16 Jun 2020 20:02
 Operator : CG/JU
 Sample : L2815-06 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OR-03-061520

Manual Integrations
 APPROVED

mohammad
 6/18/2020 11:51:38 AM

Quant Time: Jun 17 03:06:19 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 15 18:34:19 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	77442	20.00	ng	0.00
21) Naphthalene-d8	10.73	136	320248	20.00	ng	0.00
39) Acenaphthene-d10	14.57	164	212226	20.00	ng	0.00
64) Phenanthrene-d10	17.32	188	429088	20.00	ng	0.00
76) Chrysene-d12	21.58	240	397885	20.00	ng	0.00
86) Perylene-d12	24.70	264	440871	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	205644	44.86	ng	0.00
7) Phenol-d6	7.16	99	299227	42.91	ng	0.00
23) Nitrobenzene-d5	9.08	82	199288	28.43	ng	0.00
42) 2,4,6-Tribromophenol	16.06	330	116981	36.51	ng	0.00
45) 2-Fluorobiphenyl	13.19	172	454022	31.44	ng	0.00
79) Terphenyl-d14	19.93	244	638103	32.50	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.02	163	54693	3.290	ng	98
71) Phenanthrene	17.36	178	120815	5.459	ng	96
75) Fluoranthene	19.37	202	325246	11.952	ng	99
78) Pyrene	19.73	202	348576	13.083	ng	98
81) Benzo(a)anthracene	21.56	228	189129	7.332	ng	97
83) Chrysene	21.62	228	202548	8.118	ng	98
87) Indeno(1,2,3-cd)pyrene	28.25	276	156256	4.889	ng	96
88) Benzo(b)fluoranthene	23.72	252	305214	11.042	ng	97
89) Benzo(k)fluoranthene	23.77	252	90180m	3.409	ng	
90) Benzo(a)pyrene	24.55	252	217138	8.409	ng	99
92) Benzo(g,h,i)perylene	29.36	276	156674	6.107	ng	# 96

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

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