

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062620\
 Data File : BG045872.D
 Acq On : 26 Jun 2020 14:02
 Operator : CG/JU
 Sample : L3094-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BRENNAN-FL-001

Manual Integrations
 APPROVED

mohammad
 6/29/2020 3:16:58 PM

Quant Time: Jun 26 15:33:18 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 25 15:44:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	64134	20.00	ng	0.00
21) Naphthalene-d8	10.72	136	255188	20.00	ng	0.00
39) Acenaphthene-d10	14.56	164	172002	20.00	ng	0.00
64) Phenanthrene-d10	17.31	188	374599	20.00	ng	0.00
76) Chrysene-d12	21.57	240	318408	20.00	ng	0.00
86) Perylene-d12	24.70	264	358622	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	346876	91.36	ng	0.00
7) Phenol-d6	7.13	99	512373	88.72	ng	0.00
23) Nitrobenzene-d5	9.08	82	345725	61.89	ng	0.00
42) 2,4,6-Tribromophenol	16.06	330	236257	90.98	ng	0.00
45) 2-Fluorobiphenyl	13.19	172	778970	66.55	ng	0.00
79) Terphenyl-d14	19.93	244	1119117	71.23	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.01	163	137810	10.227	ng	98
71) Phenanthrene	17.36	178	109375	5.661	ng	99
75) Fluoranthene	19.37	202	119892	5.047	ng	99
78) Pyrene	19.73	202	112525	5.277	ng	97
81) Benzo(a)anthracene	21.55	228	50864m	2.464	ng	
83) Chrysene	21.61	228	46986	2.353	ng	97
88) Benzo(b)fluoranthene	23.71	252	55721	2.478	ng	93
90) Benzo(a)pyrene	24.56	252	44166	2.103	ng	97

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

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