

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP061920\
 Data File : BP002471.D
 Acq On : 19 Jun 2020 22:16
 Operator : CG/JU
 Sample : L3037-06
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SL-WC-03C

Manual Integrations
 APPROVED

mohammad
 6/22/2020 3:26:26 PM

Quant Time: Jun 20 03:49:21 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 05 15:08:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	263191	20.00	ng	0.00
21) Naphthalene-d8	10.36	136	968213	20.00	ng	0.00
39) Acenaphthene-d10	14.23	164	485771	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	889021	20.00	ng	0.00
76) Chrysene-d12	21.10	240	802837	20.00	ng	-0.01
86) Perylene-d12	23.30	264	770562	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.21	112	1904538	133.90	ng	0.00
7) Phenol-d6	6.78	99	2459029	129.85	ng	0.00
23) Nitrobenzene-d5	8.73	82	1711609	105.57	ng	0.00
42) 2,4,6-Tribromophenol	15.73	330	617960	132.87	ng	0.00
45) 2-Fluorobiphenyl	12.85	172	2957662	102.37	ng	0.00
79) Terphenyl-d14	19.58	244	3076936	105.06	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.70	163	209313	6.289	ng	Qvalue 100
71) Phenanthrene	17.03	178	273127	6.568	ng	99
72) Anthracene	17.12	178	103883	2.515	ng	99
75) Fluoranthene	19.02	202	548963	11.059	ng	99
78) Pyrene	19.37	202	543213	11.046	ng	98
81) Benzo(a)anthracene	21.08	228	298920	6.523	ng	94
83) Chrysene	21.13	228	299651	6.902	ng	98
87) Indeno(1,2,3-cd)pyrene	25.52	276	167214	3.469	ng	96
88) Benzo(b)fluoranthene	22.64	252	355855	8.564	ng	# 88
89) Benzo(k)fluoranthene	22.67	252	106722m	2.703	ng	
90) Benzo(a)pyrene	23.20	252	259131	6.654	ng	# 94
92) Benzo(a,h,i)perylene	26.19	276	199469	4.982	ng	97

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

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