

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062420\
 Data File : BP002531.D
 Acq On : 24 Jun 2020 19:51
 Operator : CG/JU
 Sample : L3096-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FK-GF-03-011-A

Manual Integrations
 APPROVED

mohammad

Quant Time: Jun 25 01:40:32 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 24 13:54:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	6/25/2020 1:07:40 PM
1) 1,4-Dichlorobenzene-d4	7.58	152	239364	20.00	ng	0.00	
21) Naphthalene-d8	10.35	136	950643	20.00	ng	0.00	
39) Acenaphthene-d10	14.23	164	537915	20.00	ng	0.00	
64) Phenanthrene-d10	16.99	188	1085271	20.00	ng	0.00	
76) Chrysene-d12	21.09	240	753962	20.00	ng	0.00	
86) Perylene-d12	23.29	264	768483	20.00	ng	0.00	
System Monitoring Compounds							
5) 2-Fluorophenol	5.20	112	2009654	155.36	ng	0.00	
7) Phenol-d6	6.78	99	2623792	152.35	ng	0.00	
23) Nitrobenzene-d5	8.73	82	1832438	115.11	ng	0.00	
42) 2,4,6-Tribromophenol	15.73	330	851651	165.36	ng	0.00	
45) 2-Fluorobiphenyl	12.85	172	3430462	107.23	ng	0.00	
79) Terphenyl-d14	19.58	244	3699901	146.08	ng	0.00	
Target Compounds							Qvalue
50) Dimethylphthalate	13.69	163	376041	10.202	ng		99
71) Phenanthrene	17.03	178	274420	5.406	ng		100
75) Fluoranthene	19.02	202	639752	10.557	ng		98
78) Pyrene	19.36	202	586400	12.697	ng		99
81) Benzo(a)anthracene	21.07	228	298569	6.938	ng		97
83) Chrysene	21.13	228	305618	7.495	ng		97
87) Indeno(1,2,3-cd)pyrene	25.50	276	208147	4.330	ng		98
88) Benzo(b)fluoranthene	22.64	252	413074	9.967	ng	#	95
89) Benzo(k)fluoranthene	22.67	252	143352m	3.641	ng		
90) Benzo(a)pyrene	23.19	252	280820	7.231	ng	#	96
92) Benzo(g,h,i)perylene	26.18	276	212650	5.325	ng		98

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

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