

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062420\
 Data File : BP002532.D
 Acq On : 24 Jun 2020 20:25
 Operator : CG/JU
 Sample : L3096-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

mohammad
 6/25/2020 1:07:43 PM

Quant Time: Jun 25 01:49:12 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)
1) 1,4-Dichlorobenzene-d4	7.58	152	247147	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	999928	20.00	ng	0.00
39) Acenaphthene-d10	14.23	164	591697	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	1180031	20.00	ng	0.00
76) Chrysene-d12	21.09	240	737745	20.00	ng	0.00
86) Perylene-d12	23.29	264	729410	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	1228024	91.94	ng	0.00
7) Phenol-d6	6.78	99	1654262	93.03	ng	0.00
23) Nitrobenzene-d5	8.73	82	1132529	67.64	ng	0.00
42) 2,4,6-Tribromophenol	15.73	330	578082	102.04	ng	0.00
45) 2-Fluorobiphenyl	12.85	172	2323844	66.04	ng	0.00
79) Terphenyl-d14	19.57	244	2600484	93.32	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.69	163	404651	9.981	ng	99
71) Phenanthrene	17.03	178	284496	5.154	ng	99
75) Fluoranthene	19.02	202	642042	9.744	ng	98
78) Pyrene	19.36	202	589618	13.048	ng	98
81) Benzo(a)anthracene	21.08	228	273223	6.488	ng	95
83) Chrysene	21.13	228	280810m	7.038	ng	
87) Indeno(1,2,3-cd)pyrene	25.50	276	190626	4.178	ng	97
88) Benzo(b)fluoranthene	22.64	252	364424m	9.265	ng	
89) Benzo(k)fluoranthene	22.67	252	127454m	3.410	ng	
90) Benzo(a)pyrene	23.20	252	256599	6.961	ng	# 95
92) Benzo(g,h,i)perylene	26.18	276	199339	5.259	ng	94

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

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