

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082219\
 Data File : BG042421.D
 Acq On : 23 Aug 2019 3:58
 Operator : HP/JU
 Sample : K4086-01RE 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NB-08-082019

Manual Integrations
 APPROVED

Jagrut
 8/23/2019 2:54:30 PM

Quant Time: Aug 23 06:49:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 22 14:29:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.00	152	68819	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	256313	20.00	ng	0.00
39) Acenaphthene-d10	14.66	164	160623	20.00	ng	0.00
64) Phenanthrene-d10	17.41	188	317744	20.00	ng	0.00
76) Chrysene-d12	21.70	240	329655	20.00	ng	0.00
87) Perylene-d12	24.94	264	358528	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	411903	121.22	ng	0.00
7) Phenol-d6	7.17	99	502792	109.54	ng	0.00
23) Nitrobenzene-d5	9.17	82	247200	66.65	ng	0.00
42) 2,4,6-Tribromophenol	16.16	330	183632	80.43	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	696220	73.31	ng	0.00
79) Terphenyl-d14	20.03	244	1007682	66.09	ng	0.00
Target Compounds						
49) Acenaphthylene	14.39	152	27880	2.070	ng	Qvalue # 94
50) Dimethylphthalate	14.11	163	136919	11.814	ng	98
71) Phenanthrene	17.46	178	109081	6.911	ng	98
75) Fluoranthene	19.47	202	371156	19.269	ng	99
78) Pyrene	19.83	202	367533	18.185	ng	99
81) Benzo(a)anthracene	21.67	228	251374	12.535	ng	98
83) Chrysene	21.74	228	278473	14.399	ng	98
86) Indeno(1,2,3-cd)pyrene	28.66	276	208060	7.916	ng	# 95
88) Benzo(b)fluoranthene	23.92	252	388913	19.731	ng	# 95
89) Benzo(k)fluoranthene	23.98	252	135024m	7.127	ng	
90) Benzo(a)pyrene	24.79	252	285920	15.287	ng	# 96
91) Dibenzo(a,h)anthracene	28.70	278	49973	2.639	ng	# 66
92) Benzo(a,h,i)perylene	29.82	276	177480	9.370	ng	95

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

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