

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082019\
 Data File : BG042367.D
 Acq On : 20 Aug 2019 21:14
 Operator : HP/JU
 Sample : K4387-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 T6-C-(0-2)

Manual Integrations
 APPROVED

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 8/22/2019 12:49:14 PM

Quant Time: Aug 21 05:27:20 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG081919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 19 13:40:37 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	78067	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	300500	20.00	ng	0.00
39) Acenaphthene-d10	14.66	164	203341	20.00	ng	0.00
64) Phenanthrene-d10	17.41	188	388703	20.00	ng	0.00
76) Chrysene-d12	21.70	240	398602	20.00	ng	0.00
87) Perylene-d12	24.94	264	478723	20.00	ng	0.00

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System Monitoring Compounds

5) 2-Fluorophenol	5.57	112	330071	85.66	ng	0.00
7) Phenol-d6	7.17	99	460385	87.98	ng	0.00
23) Nitrobenzene-d5	9.18	82	257776	60.33	ng	0.00
42) 2,4,6-Tribromophenol	16.16	330	222356	80.29	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	679362	56.54	ng	0.00
79) Terphenyl-d14	20.03	244	948999	52.15	ng	0.00

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Target Compounds

						Qvalue
15) Benzyl Alcohol	8.25	79	7248m	2.014	ng	
22) Acetophenone	8.83	105	39611	6.204	ng	# 99
50) Dimethylphthalate	14.11	163	42835	2.965	ng	99
71) Phenanthrene	17.46	178	148739	7.683	ng	99
75) Fluoranthene	19.47	202	306950	13.149	ng	100
78) Pyrene	19.83	202	298307	12.163	ng	98
81) Benzo(a)anthracene	21.67	228	181897	7.431	ng	98
83) Chrysene	21.74	228	170142	7.266	ng	97
84) Bis(2-ethylhexyl)phthalate	21.58	149	49817	3.676	ng	97
86) Indeno(1,2,3-cd)pyrene	28.63	276	102263	3.271	ng	# 95
88) Benzo(b)fluoranthene	23.91	252	212981	8.094	ng	# 97
89) Benzo(k)fluoranthene	23.97	252	71361m	2.774	ng	
90) Benzo(a)pyrene	24.78	252	156953	6.225	ng	# 98
92) Benzo(a,h,i)perylene	29.78	276	99540	3.971	ng	98

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

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