

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082219\
 Data File : BG042422.D
 Acq On : 23 Aug 2019 4:36
 Operator : HP/JU
 Sample : K4087-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 HD-02-082119

Manual Integrations
 APPROVED

Jagrut
 8/23/2019 2:54:32 PM

Quant Time: Aug 23 14:16:19 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	122640	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	478161	20.00	ng	0.00
39) Acenaphthene-d10	14.66	164	307336	20.00	ng	0.00
64) Phenanthrene-d10	17.42	188	606595	20.00	ng	0.00
76) Chrysene-d12	21.69	240	585115	20.00	ng	0.00
87) Perylene-d12	24.94	264	702973	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	672935	111.13	ng	0.00
7) Phenol-d6	7.17	99	929019	113.58	ng	0.00
23) Nitrobenzene-d5	9.17	82	528967	76.45	ng	0.00
42) 2,4,6-Tribromophenol	16.16	330	371379	85.02	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	1341177	73.81	ng	0.00
79) Terphenyl-d14	20.03	244	1764509	65.20	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.11	163	277357	12.507	ng	Qvalue 100
71) Phenanthrene	17.46	178	148243	4.920	ng	98
75) Fluoranthene	19.47	202	403639	10.977	ng	96
78) Pyrene	19.83	202	369548	10.302	ng	98
81) Benzo(a)anthracene	21.67	228	176957	4.972	ng	98
83) Chrysene	21.74	228	182942	5.329	ng	97
86) Indeno(1,2,3-cd)pyrene	28.65	276	128163	2.747	ng	# 92
88) Benzo(b)fluoranthene	23.91	252	267176	6.913	ng	# 93
89) Benzo(k)fluoranthene	23.97	252	87203m	2.347	ng	
90) Benzo(a)pyrene	24.79	252	197140	5.376	ng	# 91
92) Benzo(a,h,i)perylene	29.80	276	133188	3.586	ng	# 89

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

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