

Data Path : Z:\HPCHEM1\BNA G\DATA\BG031418\
 Data File : BG033175.D
 Acq On : 15 Mar 2018 10:46
 Operator : SJ/JU
 Sample : J1908-05DL 2X
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-BDL

Manual Integrations
 APPROVED

Sohil
 3/15/2018 7:00:26 PM

Quant Time: Mar 15 11:33:58 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 14 13:59:34 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.89	152	39931	20.00	ng	0.00
21) Naphthalene-d8	11.78	136	189344	20.00	ng	0.00
38) Acenaphthene-d10	15.51	164	118281	20.00	ng	0.00
63) Phenanthrene-d10	18.25	188	254054	20.00	ng	0.00
75) Chrysene-d12	22.60	240	237220	20.00	ng	0.00
86) Perylene-d12	26.41	264	264318	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.32	112	123974	49.67	ng	0.00
7) Phenol-d6	8.00	99	174990	50.30	ng	0.00
23) Nitrobenzene-d5	10.10	82	107195	44.62	ng	0.00
41) 2,4,6-Tribromophenol	16.99	330	69453	55.84	ng	0.00
44) 2-Fluorobiphenyl	14.14	172	248935	30.35	ng	0.00
78) Terphenyl-d14	20.76	244	338443	32.05	ng	0.00
Target Compounds						
57) Fluorene	16.55	166	29011	3.136	ng	97
70) Phenanthrene	18.29	178	491599	34.684	ng	99
71) Anthracene	18.38	178	140369	9.865	ng	99
73) Di-n-butylphthalate	19.13	149	71634	4.163	ng	99
74) Fluoranthene	20.24	202	1018849	65.074	ng	96
77) Pyrene	20.60	202	856296	51.187	ng	98
80) Benzo(a)anthracene	22.58	228	517193	34.000	ng	99
82) Chrysene	22.66	228	425787	29.302	ng	97
85) Indeno(1,2,3-cd)pyrene	30.81	276	309914	18.288	ng	98
87) Benzo(b)fluoranthene	25.18	252	609291	38.986	ng	99
88) Benzo(k)fluoranthene	25.25	252	225090m	14.492	ng	
89) Benzo(a)pyrene	26.22	252	457921	30.806	ng	99
90) Dibenzo(a,h)anthracene	30.87	278	80998m	5.413	ng	
91) Benzo(g,h,i)perylene	32.20	276	280162	18.995	ng	99

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

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