

Data Path : Z:\HPCHEM1\BNA G\DATA\BG041516\
 Data File : BG021717.D
 Acq On : 16 Apr 2016 14:40
 Operator : UM/SJ
 Sample : H2560-01
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 POSTEX-25-20160414

Manual Integrations
 APPROVED

Sohil
 4/18/2016 7:21:18 PM

Quant Time: Apr 18 03:17:10 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 13 15:48:13 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.21	152	30989	20.00	ng	0.00
21) Naphthalene-d8	11.03	136	140001	20.00	ng	0.00
38) Acenaphthene-d10	14.82	164	88233	20.00	ng	0.00
63) Phenanthrene-d10	17.56	188	195575	20.00	ng	0.00
75) Chrysene-d12	21.83	240	216376	20.00	ng	0.00
86) Perylene-d12	25.17	264	216363	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.77	112	200169	107.57	ng	0.01
7) Phenol-d6	7.37	99	305701	109.98	ng	0.01
23) Nitrobenzene-d5	9.38	82	191261	70.22	ng	0.00
41) 2,4,6-Tribromophenol	16.30	330	118516	124.63	ng	0.00
44) 2-Fluorobiphenyl	13.45	172	434476	72.15	ng	0.00
78) Terphenyl-d14	20.14	244	565666	65.23	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	14.27	163	58648	7.69	ng	100
70) Phenanthrene	17.60	178	98730	9.16	ng	96
74) Fluoranthene	19.59	202	206932	16.08	ng	100
77) Pyrene	19.95	202	238207	19.09	ng	100
80) Benzo(a)anthracene	21.81	228	126122	10.13	ng	99
82) Chrysene	21.88	228	130466	10.91	ng	97
85) Indeno(1,2,3-cd)pyrene	28.98	276	71954	5.06	ng	99
87) Benzo(b)fluoranthene	24.10	252	148743	11.85	ng	96
88) Benzo(k)fluoranthene	24.17	252	52688m	4.29	ng	
89) Benzo(a)pyrene	25.01	252	113525	9.33	ng	96
91) Benzo(g,h,i)perylene	30.16	276	78887m	6.59	ng	

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

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