

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092216\
 Data File : BG024074.D
 Acq On : 22 Sep 2016 18:27
 Operator : UM/SJ
 Sample : H4819-06DL 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YORK-SB2-07-DUPDL

Quant Time: Sep 23 01:50:52 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 20 19:01:00 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	82382	20.00	ng	0.00
21) Naphthalene-d8	10.89	136	384045	20.00	ng	0.00
38) Acenaphthene-d10	14.73	164	244575	20.00	ng	0.00
63) Phenanthrene-d10	17.49	188	509182	20.00	ng	0.00
75) Chrysene-d12	21.80	240	494917	20.00	ng	0.00
86) Perylene-d12	25.15	264	507883	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	114629	24.43	ng	0.00
7) Phenol-d6	7.23	99	180783	25.03	ng	0.00
23) Nitrobenzene-d5	9.24	82	140282	16.31	ng	0.00
41) 2,4,6-Tribromophenol	16.24	330	74587	16.93	ng	0.00
44) 2-Fluorobiphenyl	13.34	172	281510	16.50	ng	0.00
78) Terphenyl-d14	20.10	244	336139	15.47	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	14.17	163	51083	2.44	ng	97
74) Fluoranthene	19.55	202	78825	2.47	ng	99
77) Pyrene	19.91	202	68950	2.27	ng	91
83) Bis(2-ethylhexyl)phthalate	21.65	149	558115	34.74	ng	95
87) Benzo(b)fluoranthene	24.08	252	60446	2.10	ng	# 96

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

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