

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102315\
 Data File : BF082497.D
 Acq On : 24 Oct 2015 7:14
 Operator : UM/IZ
 Sample : G4116-05
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NB3-03(0-2)

Manual Integrations
 APPROVED

MmDadoda
 10/26/2015 7:22:49 PM

Quant Time: Oct 24 23:57:02 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 23 15:17:13 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	88752m	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	347865	20.00	ng	0.00
38) Acenaphthene-d10	9.79	164	166299	20.00	ng	-0.01
63) Phenanthrene-d10	11.28	188	299342	20.00	ng	-0.01
75) Chrysene-d12	13.97	240	218498	20.00	ng	-0.07
86) Perylene-d12	15.48	264	192956	20.00	ng	-0.15
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	598090	136.01	ng	0.01
7) Phenol-d6	6.41	99	802210	140.18	ng	-0.01
23) Nitrobenzene-d5	7.33	82	471975	91.11	ng	-0.01
41) 2,4,6-Tribromophenol	10.59	330	169330	108.57	ng	-0.01
44) 2-Fluorobiphenyl	9.12	172	861502	85.91	ng	-0.01
78) Terphenyl-d14	12.88	244	769331	93.30	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	9.52	163	208306	17.79	ng	99
70) Phenanthrene	11.31	178	41252	2.81	ng	97
74) Fluoranthene	12.50	202	68111	4.32	ng	98
77) Pyrene	12.73	202	59576	4.59	ng	97
80) Benzo(a)anthracene	13.96	228	30357	2.67	ng	97
82) Chrysene	13.99	228	24972	2.08	ng	98
87) Benzo(b)fluoranthene	15.08	252	33283m	2.84	ng	
89) Benzo(a)pyrene	15.42	252	22168	2.20	ng	# 97

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

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