

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100915\
 Data File : BF082157.D
 Acq On : 9 Oct 2015 21:14
 Operator : UM/IZ
 Sample : G3933-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-COMP-04

Manual Integrations
 APPROVED

MMDadoda
 10/12/2015 2:44:39 PM

Quant Time: Oct 09 23:49:37 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 09 22:47:26 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	113864	20.00	ng	-0.01
21) Naphthalene-d8	8.14	136	435646	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	179765	20.00	ng	0.00
63) Phenanthrene-d10	11.38	188	276429	20.00	ng	0.00
75) Chrysene-d12	14.02	240	241663	20.00	ng	-0.01
86) Perylene-d12	15.46	264	172418	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	582902	92.93	ng	0.01
7) Phenol-d6	6.49	99	793623	100.55	ng	0.00
23) Nitrobenzene-d5	7.42	82	431797	66.07	ng	-0.01
41) 2,4,6-Tribromophenol	10.69	330	150953	82.91	ng	-0.01
44) 2-Fluorobiphenyl	9.21	172	787369	70.98	ng	-0.01
78) Terphenyl-d14	12.97	244	611607	80.57	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.61	163	159275	12.84	ng	# 97
74) Fluoranthene	12.59	202	68424	4.62	ng	87
77) Pyrene	12.82	202	77357	5.36	ng	97
80) Benzo(a)anthracene	14.01	228	42057m	3.23	ng	
82) Chrysene	14.05	228	45939m	3.56	ng	
85) Indeno(1,2,3-cd)pyrene	16.88	276	22513	2.21	ng	94
87) Benzo(b)fluoranthene	15.05	252	47896m	4.87	ng	
88) Benzo(k)fluoranthene	15.06	252	20776m	2.30	ng	
89) Benzo(a)pyrene	15.41	252	37428	4.38	ng	# 91
91) Benzo(g,h,i)perylene	17.32	276	23154	3.15	ng	# 74

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

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