

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071015\
 Data File : BF080423.D
 Acq On : 11 Jul 2015 8:53
 Operator : TP/UM
 Sample : G2925-03 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-E4

Manual Integrations
 APPROVED

apatel
 7/13/2015 3:31:59 PM

Quant Time: Jul 13 01:46:29 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 11 05:35:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.14	152	22790	20.00	ng	-0.01
21) Naphthalene-d8	8.43	136	90126	20.00	ng	0.00
38) Acenaphthene-d10	10.19	164	44374	20.00	ng	0.00
63) Phenanthrene-d10	11.69	188	81450	20.00	ng	0.00
75) Chrysene-d12	14.57	240	96091	20.00	ng	0.03
86) Perylene-d12	16.35	264	97723	20.00	ng	0.07
System Monitoring Compounds						
5) 2-Fluorophenol	5.79	112	52468	39.45	ng	0.00
7) Phenol-d6	6.78	99	66481	37.52	ng	-0.01
23) Nitrobenzene-d5	7.71	82	38940	24.64	ng	0.00
41) 2,4,6-Tribromophenol	10.98	330	14103	32.56	ng	-0.01
44) 2-Fluorobiphenyl	9.50	172	82748	25.12	ng	0.00
78) Terphenyl-d14	13.33	244	95716	21.63	ng	0.01
Target Compounds						Qvalue
48) Acenaphthylene	10.05	152	10474	2.14	ng	97
49) Dimethylphthalate	9.90	163	7759	2.31	ng	99
70) Phenanthrene	11.71	178	54733	11.18	ng	98
71) Anthracene	11.76	178	23302	5.14	ng	99
74) Fluoranthene	12.93	202	142451	28.07	ng	93
77) Pyrene	13.19	202	148142	24.06	ng	99
80) Benzo(a)anthracene	14.56	228	96762	15.98	ng	99
82) Chrysene	14.60	228	66203	11.87	ng	97
85) Indeno(1,2,3-cd)pyrene	18.01	276	58415	8.43	ng	98
87) Benzo(b)fluoranthene	15.86	252	134284m	19.78	ng	
88) Benzo(k)fluoranthene	15.88	252	28642m	5.16	ng	
89) Benzo(a)pyrene	16.28	252	86723	14.58	ng	98
90) Dibenzo(a,h)anthracene	18.02	278	14147	2.40	ng	# 83
91) Benzo(g,h,i)perylene	18.51	276	62131	10.61	ng	99

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

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