

Data Path : V:\HPCHEM1\BNA F\DATA\BF112015\
 Data File : BF083085.D
 Acq On : 20 Nov 2015 16:52
 Operator : UM/IZ
 Sample : G4436-07DL 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ON-1593DL

Quant Time: Nov 20 17:39:32 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF111915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 20 01:33:39 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	146685	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	516676	20.00	ng	-0.01
38) Acenaphthene-d10	9.74	164	243661	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	389983	20.00	ng	0.00
75) Chrysene-d12	13.86	240	385673	20.00	ng	0.00
86) Perylene-d12	15.24	264	317070	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.30	112	637954	67.55	ng	0.01
7) Phenol-d6	6.35	99	814816	65.51	ng	0.00
23) Nitrobenzene-d5	7.26	82	524830	46.50	ng	-0.01
41) 2,4,6-Tribromophenol	10.53	330	130536	50.46	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	822538	45.46	ng	-0.01
78) Terphenyl-d14	12.81	244	552427	33.81	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
31) Naphthalene	8.01	128	95949	3.56	ng	98
37) 2-Methylnaphthalene	8.70	142	50656	2.83	ng	94
48) Acenaphthylene	9.60	152	130351	4.86	ng	98
49) Dimethylphthalate	9.47	163	226730	11.31	ng	98
51) Acenaphthene	9.77	154	35615	2.10	ng	97
54) Dibenzofuran	9.94	168	58146	2.56	ng	98
57) Fluorene	10.28	166	54707	2.99	ng	# 97
70) Phenanthrene	11.24	178	772909	32.85	ng	100
71) Anthracene	11.30	178	305410	12.89	ng	96
72) Carbazole	11.45	167	109082	5.26	ng	97
74) Fluoranthene	12.43	202	1243409	50.71	ng	99
77) Pvrene	12.66	202	1234683	45.47	ng	99
80) Benzo(a)anthracene	13.85	228	810918m	33.09	ng	
82) Chrysene	13.88	228	743860m	33.10	ng	
85) Indeno(1,2,3-cd)pyrene	16.55	276	279465	15.28	ng	95
87) Benzo(b)fluoranthene	14.87	252	844113m	38.07	ng	
88) Benzo(k)fluoranthene	14.89	252	309657m	18.33	ng	
89) Benzo(a)pyrene	15.19	252	497128	27.94	ng	# 96
90) Dibenzo(a,h)anthracene	16.56	278	91859	5.93	ng	94
91) Benzo(a,h,i)perylene	16.93	276	238846	15.81	ng	97

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

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