

Data Path : V:\HPCHEM1\BNA F\DATA\BF112015\
 Data File : BF083107.D
 Acq On : 21 Nov 2015 6:28
 Operator : UM/IZ
 Sample : G4412-08DL 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-C2DL

Quant Time: Nov 23 15:41:06 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	6.69	152	167433	20.00	ng	-0.01
21) Naphthalene-d8	7.98	136	610831	20.00	ng	-0.01
38) Acenaphthene-d10	9.74	164	295540	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	476785	20.00	ng	0.00
75) Chrysene-d12	13.85	240	385683	20.00	ng	-0.01
86) Perylene-d12	15.23	264	340167	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	694121	64.39	ng	0.01
7) Phenol-d6	6.35	99	870043	61.28	ng	0.00
23) Nitrobenzene-d5	7.26	82	605785	45.40	ng	-0.01
41) 2,4,6-Tribromophenol	10.53	330	135915	43.31	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	924397	42.13	ng	-0.01
78) Terphenyl-d14	12.81	244	641581	39.27	ng	0.00

Target Compounds				Qvalue		
31) Naphthalene	8.01	128	232995	7.32	ng	99
37) 2-Methylnaphthalene	8.70	142	76914	3.63	ng	94
48) Acenaphthylene	9.60	152	189801	5.83	ng	99
49) Dimethylphthalate	9.47	163	240515	9.89	ng	97
54) Dibenzofuran	9.94	168	123486	4.48	ng	99
57) Fluorene	10.28	166	113117	5.10	ng	99
70) Phenanthrene	11.24	178	1261920	43.87	ng	98
71) Anthracene	11.29	178	133073	4.59	ng	98
72) Carbazole	11.45	167	102139	4.03	ng	99
74) Fluoranthene	12.43	202	1282953	42.79	ng	99
77) Pyrene	12.66	202	1352992	49.83	ng	99
80) Benzo(a)anthracene	13.85	228	582746m	23.78	ng	
82) Chrysene	13.88	228	589828	26.24	ng	98
83) Bis(2-ethylhexyl)phthalate	13.86	149	34314	2.19	ng	# 95
85) Indeno(1,2,3-cd)pyrene	16.53	276	246002	13.45	ng	98
87) Benzo(b)fluoranthene	14.85	252	608446m	25.58	ng	
88) Benzo(k)fluoranthene	14.87	252	242399m	13.38	ng	
89) Benzo(a)pyrene	15.19	252	423092	22.17	ng	98
90) Dibenzo(a,h)anthracene	16.53	278	57255m	3.45	ng	
91) Benzo(a,h,i)perylene	16.92	276	225166	13.90	ng	98

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

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