

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112415\
 Data File : BF083238.D
 Acq On : 24 Nov 2015 15:50
 Operator : UM/IZ
 Sample : G4503-07
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-7

Quant Time: Nov 24 17:09:41 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 24 15:18:40 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.60	152	174700	20.00	ng	-0.01
21) Naphthalene-d8	7.90	136	660726	20.00	ng	-0.01
38) Acenaphthene-d10	9.64	164	329640	20.00	ng	-0.01
63) Phenanthrene-d10	11.13	188	621831	20.00	ng	0.00
75) Chrysene-d12	13.75	240	487116	20.00	ng	-0.01
86) Perylene-d12	15.11	264	373964	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.21	112	738367	63.90	ng	0.01
7) Phenol-d6	6.28	99	631000	42.20	ng	0.00
23) Nitrobenzene-d5	7.19	82	1498332	103.62	ng	0.00
41) 2,4,6-Tribromophenol	10.44	330	473717	140.47	ng	0.00
44) 2-Fluorobiphenyl	8.98	172	2026293	85.72	ng	-0.01
78) Terphenyl-d14	12.72	244	1906393	92.16	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	6.30	94	55307	3.06	ng	81
17) 2-Methylphenol	6.90	107	42745	4.14	ng	# 83
20) 3+4-Methylphenols	7.06	107	147353	11.28	ng	95
22) Acetophenone	7.03	105	54420	2.87	ng	# 94
27) 2,4-Dimethylphenol	7.59	122	59321	5.45	ng	90
31) Naphthalene	7.92	128	1604199	44.99	ng	98
37) 2-Methylnaphthalene	8.62	142	184260	7.96	ng	96
49) Dimethylphthalate	9.38	163	70410	2.77	ng	98
51) Acenaphthene	9.68	154	203000	9.77	ng	97
54) Dibenzofuran	9.85	168	208691	7.10	ng	97
57) Fluorene	10.19	166	198737	8.18	ng	99
70) Phenanthrene	11.15	178	591030	16.50	ng	98
71) Anthracene	11.20	178	111261	3.04	ng	97
72) Carbazole	11.37	167	412008	12.68	ng	97
74) Fluoranthene	12.33	202	117038	3.19	ng	99
77) Pyrene	12.56	202	79716	2.40	ng	97

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

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