

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022315\
 Data File : BF077403.D
 Acq On : 23 Feb 2015 14:16
 Operator : TP/IZ
 Sample : G1332-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HF-3

Quant Time: Feb 24 02:26:24 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 23 23:58:25 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.33	152	76212	20.00	ng	0.00
21) Naphthalene-d8	8.91	136	312086	20.00	ng	0.00
38) Acenaphthene-d10	11.08	164	171992	20.00	ng	-0.01
63) Phenanthrene-d10	12.95	188	296282	20.00	ng	0.01
75) Chrysene-d12	16.24	240	316667	20.00	ng	-0.02
86) Perylene-d12	18.01	264	352797	20.00	ng	-0.09

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.62	112	447778	98.09	ng	0.00
7) Phenol-d6	6.88	99	569249	96.98	ng	-0.01
23) Nitrobenzene-d5	8.02	82	351671	70.07	ng	-0.01
41) 2,4,6-Tribromophenol	12.08	330	134399	81.16	ng	0.01
44) 2-Fluorobiphenyl	10.26	172	687455	62.32	ng	0.00
78) Terphenyl-d14	14.93	244	765707	56.53	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	6.89	94	14083	2.02	ng	# 30
31) Naphthalene	8.94	128	5150000	329.98	ng	# 90
37) 2-Methylnaphthalene	9.80	142	3359018	303.07	ng	# 95
49) Dimethylphthalate	10.76	163	47936	3.97	ng	# 97
51) Acenaphthene	11.14	154	3203481	331.83	ng	97
54) Dibenzofuran	11.36	168	4085207	274.07	ng	# 87
57) Fluorene	11.79	166	4624576	388.07	ng	97
70) Phenanthrene	13.00	178	11193469	651.83	ng	# 85
71) Anthracene	13.05	178	2396839	140.65	ng	95
72) Carbazole	13.23	167	826044	50.98	ng	97
74) Fluoranthene	14.48	202	6506294	346.85	ng	90
77) Pvrene	14.75	202	5651394	291.75	ng	# 90
80) Benzo(a)anthracene	16.23	228	2079377	112.04	ng	94
82) Chrysene	16.27	228	1526538	87.60	ng	96
85) Indeno(1,2,3-cd)pyrene	19.49	276	210842m	10.61	ng	
87) Benzo(b)fluoranthene	17.55	252	984868m	48.01	ng	
88) Benzo(k)fluoranthene	17.57	252	399922m	19.89	ng	
89) Benzo(a)pyrene	17.94	252	649100	33.22	ng	100
90) Dibenzo(a,h)anthracene	19.53	278	59790	3.21	ng	# 93
91) Benzo(a,h,i)perylene	19.94	276	180327	9.59	ng	99

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

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