

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011915\
 Data File : BF076952.D
 Acq On : 20 Jan 2015 6:19
 Operator : IZ / TP
 Sample : G1076-01
 Misc : W
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-1

Quant Time: Jan 20 10:31:11 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 20 06:05:35 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.98	152	37629	20.00	ng	0.01
21) Naphthalene-d8	10.88	136	152117	20.00	ng	0.00
38) Acenaphthene-d10	15.01	164	81929	20.00	ng	0.00
63) Phenanthrene-d10	17.79	188	132040	20.00	ng	0.01
75) Chrysene-d12	21.28	240	121505	20.00	ng	0.01
86) Perylene-d12	22.92	264	118010	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.04	112	203130	88.76	ng	0.00
7) Phenol-d6	7.37	99	293339	101.60	ng	-0.01
23) Nitrobenzene-d5	9.27	82	182858	66.60	ng	0.00
41) 2,4,6-Tribromophenol	16.68	330	40891	61.49	ng	-0.01
44) 2-Fluorobiphenyl	13.53	172	336731	62.55	ng	0.00
78) Terphenyl-d14	20.00	244	302126	57.24	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	7.41	94	11415	3.72	ng	# 66
31) Naphthalene	10.93	128	17785	2.27	ng	98
37) 2-Methylnaphthalene	12.59	142	21589	4.04	ng	94
49) Dimethylphthalate	14.60	163	45047	7.75	ng	# 98
51) Acenaphthene	15.09	154	95543	19.97	ng	96
54) Dibenzofuran	15.51	168	77909	11.18	ng	97
57) Fluorene	16.23	166	91252	15.46	ng	98
59) Diethylphthalate	16.21	149	52550	8.95	ng	97
70) Phenanthrene	17.82	178	1113395	135.22	ng	99
71) Anthracene	17.89	178	282743	35.21	ng	99
72) Carbazole	18.17	167	84484	10.85	ng	98
74) Fluoranthene	19.47	202	1685763	199.80	ng	97
77) Pvrene	19.75	202	1361610	163.51	ng	# 89
80) Benzo(a)anthracene	21.27	228	787520	103.31	ng	99
82) Chrysene	21.32	228	626816m	90.16	ng	
85) Indeno(1,2,3-cd)pyrene	24.03	276	394707m	53.57	ng	
87) Benzo(b)fluoranthene	22.53	252	873034m	109.84	ng	
88) Benzo(k)fluoranthene	22.55	252	244805m	35.25	ng	
89) Benzo(a)pvrene	22.87	252	581896m	83.00	ng	
90) Dibenzo(a,h)anthracene	24.04	278	77222m	12.00	ng	
91) Benzo(q,h,i)perylene	24.32	276	341232m	53.36	ng	

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

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