

Data Path : Z:\HPCHEM1\BNA F\DATA\BF021318\
 Data File : BF102962.D
 Acq On : 13 Feb 2018 15:30
 Operator : SJ/JU
 Sample : J1511-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MM-2

Manual Integrations
 APPROVED

Sohil
 2/14/2018 10:29:38 AM

Quant Time: Feb 14 01:15:03 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 05 00:38:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	168610	20.00	ng	-0.01
21) Naphthalene-d8	8.16	136	710811	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	304931	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	445317	20.00	ng	0.00
75) Chrysene-d12	14.04	240	289687	20.00	ng	0.00
86) Perylene-d12	15.52	264	307443	20.00	ng	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.49	112	661927	59.62	ng	0.00
7) Phenol-d6	6.50	99	1110862	83.10	ng	0.00
23) Nitrobenzene-d5	7.43	82	736966	71.92	ng	-0.01
41) 2,4,6-Tribromophenol	10.70	330	59716	24.33	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	917645	49.94	ng	-0.01
78) Terphenyl-d14	12.98	244	539553	44.10	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	6.51	94	71466	4.988	ng	86
31) Naphthalene	8.17	128	816217	23.503	ng	99
37) 2-Methylnaphthalene	8.87	142	215932	9.497	ng	100
45) 1,1'-Biphenyl	9.33	154	65796	2.562	ng	98
49) Dimethylphthalate	9.62	163	109940	4.624	ng	# 96
51) Acenaphthene	9.95	154	170184	9.061	ng	99
54) Dibenzofuran	10.12	168	254746	9.625	ng	96
57) Fluorene	10.46	166	176884	9.185	ng	99
65) n-Nitrosodiphenylamine	10.56	169	36436	2.320	ng	# 57
70) Phenanthrene	11.43	178	1744579	70.342	ng	99
71) Anthracene	11.47	178	451064	17.881	ng	98
72) Carbazole	11.63	167	191298	7.741	ng	99
73) Di-n-butylphthalate	11.95	149	85588	2.851	ng	98
74) Fluoranthene	12.62	202	1099252	44.053	ng	99
77) Pyrene	12.84	202	1124867	49.519	ng	100
79) Butylbenzylphthalate	13.45	149	15194	2.393	ng	# 79
80) Benzo(a)anthracene	14.03	228	514302	28.230	ng	99
82) Chrysene	14.06	228	474550	25.840	ng	98
83) Bis(2-ethylhexyl)phthalate	14.00	149	75707	5.440	ng	98
85) Indeno(1,2,3-cd)pyrene	17.00	276	199493	13.097	ng	96
87) Benzo(b)fluoranthene	15.09	252	464309m	23.294	ng	
88) Benzo(k)fluoranthene	15.11	252	184940m	9.710	ng	
89) Benzo(a)pyrene	15.45	252	379831	21.170	ng	# 94
90) Dibenzo(a,h)anthracene	17.01	278	54120	3.716	ng	# 86
91) Benzo(g,h,i)perylene	17.46	276	214687	15.062	ng	95

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

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