

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022118\
 Data File : BF103131.D
 Acq On : 21 Feb 2018 12:57
 Operator : SJ/JU
 Sample : J1403-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-04-022018

Manual Integrations
 APPROVED

Sohil
 2/22/2018 9:04:58 AM

Quant Time: Feb 21 13:43:17 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 16 14:34:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	148055	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	595462	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	237687	20.00	ng	-0.01
63) Phenanthrene-d10	11.40	188	344350	20.00	ng	0.00
75) Chrysene-d12	14.05	240	275230	20.00	ng	0.00
86) Perylene-d12	15.54	264	278153	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	947628	97.20	ng	0.00
7) Phenol-d6	6.51	99	1114893	94.98	ng	0.00
23) Nitrobenzene-d5	7.44	82	678090	79.00	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	164763	86.12	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	838433	58.53	ng	-0.01
78) Terphenyl-d14	12.99	244	615278	52.93	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.18	128	689063	23.685	ng	99
37) 2-Methylnaphthalene	8.87	142	311897	16.375	ng	98
45) 1,1'-Biphenyl	9.33	154	64568	3.225	ng	93
48) Acenaphthylene	9.77	152	254020	10.377	ng	99
49) Dimethylphthalate	9.62	163	75047	4.049	ng	# 99
51) Acenaphthene	9.95	154	121410	8.293	ng	99
54) Dibenzofuran	10.12	168	256829	12.449	ng	97
57) Fluorene	10.46	166	340067	22.655	ng	99
70) Phenanthrene	11.43	178	1261213	65.763	ng	99
71) Anthracene	11.48	178	418234	21.441	ng	98
72) Carbazole	11.63	167	129653	6.785	ng	99
74) Fluoranthene	12.63	202	1078869	55.913	ng	99
77) Pvrene	12.86	202	959836	44.473	ng	99
80) Benzo(a)anthracene	14.04	228	550540	31.806	ng	98
82) Chrysene	14.07	228	455917	26.129	ng	97
85) Indeno(1,2,3-cd)pyrene	17.06	276	194692	13.453	ng	96
87) Benzo(b)fluoranthene	15.11	252	600275m	33.286	ng	
88) Benzo(k)fluoranthene	15.13	252	179652m	10.426	ng	
89) Benzo(a)pvrene	15.48	252	391467	24.116	ng	# 94
90) Dibenzo(a,h)anthracene	17.07	278	51057	3.875	ng	# 84
91) Benzo(g,h,i)perylene	17.52	276	170900	13.252	ng	95

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

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