

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061218\
 Data File : BF106362.D
 Acq On : 12 Jun 2018 23:11
 Operator : JU/SJ
 Sample : J3322-01
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB02-2_E_5

Manual Integrations
 APPROVED

Sohil
 6/13/2018 4:56:24 PM

Quant Time: Jun 13 04:12:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 11 15:00:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	139325	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	479215	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	204875	20.00	ng	0.00
63) Phenanthrene-d10	11.51	188	420937	20.00	ng	0.00
75) Chrysene-d12	14.17	240	387207	20.00	ng	0.00
86) Perylene-d12	15.71	264	287294	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	795608	96.65	ng	0.04
7) Phenol-d6	6.63	99	1002981	96.44	ng	0.04
23) Nitrobenzene-d5	7.54	82	562486	69.05	ng	0.00
41) 2,4,6-Tribromophenol	10.81	330	217435	110.30	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	878983	72.67	ng	0.00
78) Terphenyl-d14	13.10	244	1007522	64.63	ng	0.01
Target Compounds						
31) Naphthalene	8.29	128	3065398	141.503	ng	Qvalue # 87
37) 2-Methylnaphthalene	8.97	142	146628	9.938	ng	98
45) 1,1'-Biphenyl	9.43	154	169515	10.682	ng	98
48) Acenaphthylene	9.88	152	208161	10.709	ng	99
49) Dimethylphthalate	9.72	163	108729	7.428	ng	100
51) Acenaphthene	10.05	154	367721	29.294	ng	99
54) Dibenzofuran	10.22	168	317996	18.812	ng	98
57) Fluorene	10.57	166	384119	28.681	ng	99
70) Phenanthrene	11.54	178	1307232	63.412	ng	96
71) Anthracene	11.59	178	465946	22.561	ng	99
72) Carbazole	11.75	167	320402	16.804	ng	98
74) Fluoranthene	12.75	202	1222525	55.608	ng	98
77) Pvrene	12.98	202	1063504	43.871	ng	97
80) Benzo(a)anthracene	14.15	228	624101m	27.085	ng	
82) Chrysene	14.19	228	554708m	26.367	ng	
85) Indeno(1,2,3-cd)pyrene	17.27	276	215535	12.837	ng	# 88
87) Benzo(b)fluoranthene	15.25	252	585379m	33.707	ng	
88) Benzo(k)fluoranthene	15.28	252	150233m	8.735	ng	
89) Benzo(a)pvrene	15.64	252	383835	24.572	ng	96
90) Dibenzo(a,h)anthracene	17.28	278	59593m	4.576	ng	
91) Benzo(q,h,i)perylene	17.75	276	194386	15.359	ng	94

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

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