

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062118\
 Data File : BF106689.D
 Acq On : 22 Jun 2018 4:01
 Operator : JU/SJ
 Sample : J3603-03
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 26KV-TP-2

Manual Integrations
 APPROVED

Sohil
 6/22/2018 10:23:48 AM

Quant Time: Jun 22 04:47:00 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 19 14:32:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	58382	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	205458	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	93071	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	179691	20.00	ng	0.00
75) Chrysene-d12	14.15	240	142711	20.00	ng	0.00
86) Perylene-d12	15.68	264	107731	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	332099	96.32	ng	0.00
7) Phenol-d6	6.61	99	406223	94.35	ng	0.00
23) Nitrobenzene-d5	7.52	82	242165	65.50	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	96945	89.87	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	335032	55.88	ng	0.00
78) Terphenyl-d14	13.08	244	359346	60.99	ng	0.00
Target Compounds						
31) Naphthalene	8.27	128	32395	3.372	ng	98
49) Dimethylphthalate	9.71	163	45268	6.485	ng	98
51) Acenaphthene	10.04	154	26332	4.536	ng	99
54) Dibenzofuran	10.21	168	18011	2.266	ng	98
57) Fluorene	10.55	166	30231	4.793	ng	97
70) Phenanthrene	11.52	178	397152	45.824	ng	99
71) Anthracene	11.57	178	93377	10.555	ng	99
72) Carbazole	11.72	167	37042	4.472	ng	97
74) Fluoranthene	12.72	202	599613	62.769	ng	97
77) Pyrene	12.95	202	634737	67.448	ng	99
80) Benzo(a)anthracene	14.14	228	297981	34.259	ng	96
82) Chrysene	14.17	228	270225	33.770	ng	98
85) Indeno(1,2,3-cd)pyrene	17.25	276	88472	13.028	ng	# 88
87) Benzo(b)fluoranthene	15.22	252	257347m	38.455	ng	
88) Benzo(k)fluoranthene	15.25	252	59327m	9.309	ng	
89) Benzo(a)pyrene	15.62	252	176966m	29.746	ng	
90) Dibenzo(a,h)anthracene	17.25	278	25168m	4.992	ng	
91) Benzo(g,h,i)perylene	17.72	276	86275	17.507	ng	# 90

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

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