

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062618\
 Data File : BF106822.D
 Acq On : 26 Jun 2018 15:01
 Operator : JU/SJ
 Sample : J3676-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-1-4

Manual Integrations
 APPROVED

Sohil
 6/27/2018 2:14:02 PM

Quant Time: Jun 27 09:09:18 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 25 18:50:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	48095	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	181922	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	84466	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	143268	20.00	ng	0.00
75) Chrysene-d12	14.16	240	129046	20.00	ng	0.00
86) Perylene-d12	15.72	264	100068	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	292255	102.90	ng	0.00
7) Phenol-d6	6.61	99	365610	103.08	ng	0.00
23) Nitrobenzene-d5	7.52	82	228816	69.90	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	102467	104.67	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	345429	65.91	ng	0.00
78) Terphenyl-d14	13.08	244	323805	60.78	ng	0.00
Target Compounds						
31) Naphthalene	8.27	128	54325	6.387	ng	98
37) 2-Methylnaphthalene	8.95	142	17620	3.125	ng	92
49) Dimethylphthalate	9.71	163	89343	14.103	ng	99
51) Acenaphthene	10.04	154	21064	3.999	ng	97
54) Dibenzofuran	10.21	168	36185	5.016	ng	94
57) Fluorene	10.55	166	46634	8.147	ng	99
70) Phenanthrene	11.52	178	316323	45.777	ng	99
71) Anthracene	11.57	178	88567	12.557	ng	99
72) Carbazole	11.73	167	23358	3.537	ng	98
74) Fluoranthene	12.72	202	356897	46.859	ng	95
77) Pyrene	12.95	202	308566	36.261	ng	100
80) Benzo(a)anthracene	14.15	228	169336	21.530	ng	99
82) Chrysene	14.19	228	145346	20.087	ng	98
83) Bis(2-ethylhexyl)phthalate	14.11	149	10742	2.401	ng	# 90
85) Indeno(1,2,3-cd)pyrene	17.30	276	51565	8.398	ng	# 88
87) Benzo(b)fluoranthene	15.26	252	160483m	25.817	ng	
88) Benzo(k)fluoranthene	15.29	252	40718m	6.878	ng	
89) Benzo(a)pyrene	15.59	252	67971	12.300	ng	98
90) Dibenzo(a,h)anthracene	17.31	278	13780	2.942	ng	# 81
91) Benzo(a,h,i)perylene	17.79	276	49506	10.815	ng	# 88

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

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