

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061218\
 Data File : BF106346.D
 Acq On : 12 Jun 2018 16:00
 Operator : JU/SJ
 Sample : J3245-07 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-061118-D

Manual Integrations
 APPROVED

Sohil
 6/13/2018 4:56:09 PM

Quant Time: Jun 13 02:06:11 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	6.97	152	170419	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	589571	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	248275	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	484190	20.00	ng	0.00
75) Chrysene-d12	14.16	240	518601	20.00	ng	0.00
86) Perylene-d12	15.71	264	384040	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	542381	53.87	ng	0.00
7) Phenol-d6	6.59	99	652934	51.33	ng	0.00
23) Nitrobenzene-d5	7.53	82	405527	40.46	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	145505	60.91	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	698924	40.87	ng	0.00
78) Terphenyl-d14	13.09	244	846487	40.54	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.72	163	117103	6.602	ng	# 99
57) Fluorene	10.56	166	46033	2.836	ng	# 97
70) Phenanthrene	11.53	178	412056	17.377	ng	98
71) Anthracene	11.58	178	118356	4.982	ng	96
74) Fluoranthene	12.72	202	652198	25.791	ng	99
77) Pyrene	12.95	202	589844	18.167	ng	100
80) Benzo(a)anthracene	14.15	228	309192	10.019	ng	96
82) Chrysene	14.19	228	293015	10.399	ng	98
85) Indeno(1,2,3-cd)pyrene	17.28	276	87714m	3.901	ng	
87) Benzo(b)fluoranthene	15.25	252	290436m	12.511	ng	
88) Benzo(k)fluoranthene	15.28	252	72896m	3.171	ng	
89) Benzo(a)pyrene	15.65	252	184581	8.840	ng	96
91) Benzo(a,h,i)perylene	17.75	276	84864	5.016	ng	# 91

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

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