

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061518\
 Data File : BF106500.D
 Acq On : 15 Jun 2018 19:00
 Operator : JU/SJ
 Sample : J3486-06
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DOCUMENTATION-9

Manual Integrations
 APPROVED

Sohil
 6/18/2018 10:38:28 AM

Quant Time: Jun 16 00:36:25 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 15 11:11:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	137238	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	502600	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	214086	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	344579	20.00	ng	0.00
75) Chrysene-d12	14.15	240	345969	20.00	ng	0.00
86) Perylene-d12	15.69	264	303872	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	708021	87.32	ng	0.00
7) Phenol-d6	6.61	99	856741	83.63	ng	0.00
23) Nitrobenzene-d5	7.52	82	574744	67.27	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	182171	88.44	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	857601	66.04	ng	0.00
78) Terphenyl-d14	13.08	244	710413	51.00	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	134275	8.779	ng	99
70) Phenanthrene	11.52	178	235481	13.954	ng	98
71) Anthracene	11.57	178	62981	3.725	ng	98
74) Fluoranthene	12.72	202	538453	29.920	ng	96
77) Pyrene	12.95	202	498628	23.021	ng	99
80) Benzo(a)anthracene	14.14	228	370990m	18.019	ng	
82) Chrysene	14.18	228	308233	16.398	ng	97
85) Indeno(1,2,3-cd)pyrene	17.25	276	111231	7.414	ng	# 89
87) Benzo(b)fluoranthene	15.24	252	404445m	22.018	ng	
88) Benzo(k)fluoranthene	15.27	252	100216m	5.509	ng	
89) Benzo(a)pyrene	15.62	252	262399	15.882	ng	97
90) Dibenzo(a,h)anthracene	17.27	278	28637	2.079	ng	97
91) Benzo(a,h,i)perylene	17.73	276	101115	7.554	ng	94

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

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