

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
 Data File : BF106742.D
 Acq On : 23 Jun 2018 6:53
 Operator : JU/SJ
 Sample : J3581-07 2X
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSP-C-COMPOSITE

Manual Integrations
 APPROVED

Sohil
 6/25/2018 5:17:15 PM

Quant Time: Jun 25 02:26:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 22 11:34:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	51826	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	187335	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	78911	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	146957	20.00	ng	-0.01
75) Chrysene-d12	14.14	240	148706	20.00	ng	-0.01
86) Perylene-d12	15.68	264	109159	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	179975	58.80	ng	0.00
7) Phenol-d6	6.60	99	222897	58.32	ng	0.00
23) Nitrobenzene-d5	7.52	82	137074	40.66	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	54666	59.77	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	255599	48.49	ng	-0.01
78) Terphenyl-d14	13.08	244	285289	46.47	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	31333	5.294	ng	98
70) Phenanthrene	11.51	178	31584	4.456	ng	99
74) Fluoranthene	12.71	202	71457	9.147	ng	97
77) Pyrene	12.94	202	71757	7.318	ng	99
80) Benzo(a)anthracene	14.13	228	40214	4.437	ng	96
82) Chrysene	14.17	228	36026	4.321	ng	95
83) Bis(2-ethylhexyl)phthalate	14.10	149	10831	2.101	ng	# 95
85) Indeno(1,2,3-cd)pyrene	17.24	276	17088	2.415	ng	# 84
87) Benzo(b)fluoranthene	15.22	252	38269m	5.644	ng	
89) Benzo(a)pyrene	15.61	252	24155	4.007	ng	93
91) Benzo(g,h,i)perylene	17.72	276	16259	3.256	ng	# 90

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

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