

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
 Data File : BF106741.D
 Acq On : 23 Jun 2018 6:27
 Operator : JU/SJ
 Sample : J3581-01 2X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSP-A-COMPOSITE

Manual Integrations
 APPROVED

Sohil
 6/25/2018 5:17:13 PM

Quant Time: Jun 25 02:18:10 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	49824	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	182169	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	78574	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	141639	20.00	ng	-0.01
75) Chrysene-d12	14.14	240	149721	20.00	ng	-0.01
86) Perylene-d12	15.68	264	111425	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	131160	44.58	ng	0.00
7) Phenol-d6	6.60	99	163348	44.45	ng	0.00
23) Nitrobenzene-d5	7.52	82	99454	30.34	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	39492	43.36	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	192422	32.31	ng	-0.01
78) Terphenyl-d14	13.07	244	209880	33.96	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.71	163	20664	3.506	ng	99
70) Phenanthrene	11.51	178	23876	3.495	ng	100
74) Fluoranthene	12.71	202	51323	6.816	ng	95
77) Pyrene	12.94	202	52970	5.365	ng	99
80) Benzo(a)anthracene	14.13	228	29456m	3.228	ng	
82) Chrysene	14.17	228	26167m	3.117	ng	
87) Benzo(b)fluoranthene	15.22	252	32872m	4.749	ng	
89) Benzo(a)pyrene	15.61	252	20134	3.272	ng	# 91
91) Benzo(g,h,i)perylene	17.72	276	12364	2.426	ng	91

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

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