

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
 Data File : BF106743.D
 Acq On : 23 Jun 2018 7:20
 Operator : JU/SJ
 Sample : J3581-04 2X
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSP-B-COMPOSITE

Manual Integrations
 APPROVED

Sohil
 6/25/2018 5:16:23 PM

Quant Time: Jun 25 07:31:29 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	49681	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	179241	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	78871	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	144216	20.00	ng	-0.01
75) Chrysene-d12	14.14	240	137973	20.00	ng	-0.01
86) Perylene-d12	15.68	264	96661	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	168699	57.50	ng	0.00
7) Phenol-d6	6.60	99	204844	55.91	ng	0.00
23) Nitrobenzene-d5	7.52	82	124998	38.76	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	51416	56.25	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	236554	43.58	ng	-0.01
78) Terphenyl-d14	13.08	244	264398	46.42	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	26087	4.410	ng	99
70) Phenanthrene	11.52	178	19698	2.832	ng	99
74) Fluoranthene	12.71	202	46309	6.040	ng	97
77) Pyrene	12.94	202	47414	5.211	ng	97
80) Benzo(a)anthracene	14.13	228	25417	3.023	ng	# 94
82) Chrysene	14.17	228	23132m	2.990	ng	
87) Benzo(b)fluoranthene	15.22	252	25571m	4.259	ng	
89) Benzo(a)pyrene	15.61	252	16472	3.086	ng	# 85
91) Benzo(g,h,i)perylene	17.73	276	12988	2.937	ng	# 92

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

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