

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061418\
 Data File : BF106459.D
 Acq On : 14 Jun 2018 22:51
 Operator : JU/SJ
 Sample : J3471-02
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 #1

Manual Integrations
 APPROVED

Sohil
 6/15/2018 11:16:55 AM

Quant Time: Jun 15 04:00:40 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 14 18:39:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	117427	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	422935	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	171641	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	301085	20.00	ng	0.00
75) Chrysene-d12	14.15	240	284803	20.00	ng	-0.01
86) Perylene-d12	15.68	264	196648	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	698385	100.66	ng	0.00
7) Phenol-d6	6.61	99	847241	96.65	ng	0.00
23) Nitrobenzene-d5	7.53	82	536198	74.58	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	177562	107.52	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	813146	83.74	ng	0.00
78) Terphenyl-d14	13.08	244	769687	67.12	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	95679	7.802	ng	99
70) Phenanthrene	11.52	178	41540	2.817	ng	97
74) Fluoranthene	12.71	202	109941	6.991	ng	98
77) Pyrene	12.95	202	107144	6.009	ng	97
80) Benzo(a)anthracene	14.14	228	54336	3.206	ng	97
82) Chrysene	14.17	228	52215	3.374	ng	99
83) Bis(2-ethylhexyl)phthalate	14.11	149	90090	9.414	ng	96
87) Benzo(b)fluoranthene	15.22	252	55606m	4.678	ng	
89) Benzo(a)pyrene	15.61	252	34607	3.237	ng	# 86
91) Benzo(g,h,i)perylene	17.72	276	23461	2.708	ng	96

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

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