

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061418\
 Data File : BF106460.D
 Acq On : 14 Jun 2018 23:18
 Operator : JU/SJ
 Sample : J3471-04
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 #2

Manual Integrations
 APPROVED

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

Quant Time: Jun 15 04:05:45 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	125973	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	451319	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	184991	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	326699	20.00	ng	0.00
75) Chrysene-d12	14.15	240	291548	20.00	ng	-0.01
86) Perylene-d12	15.68	264	205437	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	675275	90.73	ng	0.00
7) Phenol-d6	6.61	99	815950	86.77	ng	0.00
23) Nitrobenzene-d5	7.52	82	519213	67.68	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	175041	98.34	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	798020	73.23	ng	0.00
78) Terphenyl-d14	13.08	244	753943	64.23	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	86646	6.556	ng	99
70) Phenanthrene	11.52	178	55306	3.457	ng	97
74) Fluoranthene	12.71	202	138242	8.102	ng	98
77) Pyrene	12.94	202	136607	7.484	ng	99
80) Benzo(a)anthracene	14.14	228	67287	3.878	ng	98
82) Chrysene	14.17	228	64457	4.069	ng	97
87) Benzo(b)fluoranthene	15.22	252	66929m	5.389	ng	
89) Benzo(a)pyrene	15.61	252	42870	3.838	ng	# 90
91) Benzo(g,h,i)perylene	17.72	276	25688	2.838	ng	# 91

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

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