

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061418\
 Data File : BF106466.D
 Acq On : 15 Jun 2018 2:25
 Operator : JU/SJ
 Sample : J3481-03
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-021-B

Manual Integrations
 APPROVED

Sohil
 6/15/2018 11:16:57 AM

Quant Time: Jun 15 08:33:54 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	97834	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	322210	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	131976	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	292663	20.00	ng	0.00
75) Chrysene-d12	14.15	240	304079	20.00	ng	-0.01
86) Perylene-d12	15.68	264	216050	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	725924	125.58	ng	0.00
7) Phenol-d6	6.61	99	843140	115.45	ng	0.00
23) Nitrobenzene-d5	7.53	82	521801	95.27	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	210522	165.79	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	756469	112.24	ng	0.00
78) Terphenyl-d14	13.08	244	1005828	82.16	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	80737	8.563	ng	Qvalue # 99
70) Phenanthrene	11.52	178	78204	5.456	ng	99
74) Fluoranthene	12.71	202	135191	8.845	ng	96
77) Pyrene	12.94	202	135081	7.096	ng	99
80) Benzo(a)anthracene	14.14	228	67247	3.716	ng	96
82) Chrysene	14.17	228	64581	3.909	ng	96
87) Benzo(b)fluoranthene	15.22	252	57487m	4.402	ng	
89) Benzo(a)pyrene	15.61	252	36600	3.116	ng	93
91) Benzo(g,h,i)perylene	17.72	276	21716	2.282	ng	98

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

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