

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071318\
 Data File : BF107370.D
 Acq On : 13 Jul 2018 16:35
 Operator : JU/SJ
 Sample : J3825-07 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-071118-D

Manual Integrations
 APPROVED

Sohil
 7/16/2018 3:22:27 PM

Quant Time: Jul 16 11:04:04 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 13 11:54:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	46628	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	167344	20.00	ng	0.00
38) Acenaphthene-d10	9.98	164	77931	20.00	ng	0.00
63) Phenanthrene-d10	11.48	188	163089	20.00	ng	0.00
75) Chrysene-d12	14.14	240	152350	20.00	ng	0.00
86) Perylene-d12	15.68	264	108046	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	143689	52.18	ng	0.00
7) Phenol-d6	6.58	99	168507	49.00	ng	0.00
23) Nitrobenzene-d5	7.50	82	96318	31.99	ng	0.00
41) 2,4,6-Tribromophenol	10.78	330	44362	49.11	ng	0.00
44) 2-Fluorobiphenyl	9.29	172	191603	32.51	ng	0.00
78) Terphenyl-d14	13.06	244	263103	41.83	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.68	163	30398	5.201	ng	Qvalue # 98
70) Phenanthrene	11.50	178	35075	4.459	ng	97
74) Fluoranthene	12.69	202	92132	10.626	ng	96
77) Pyrene	12.93	202	96227	9.578	ng	98
80) Benzo(a)anthracene	14.12	228	40311	4.341	ng	98
82) Chrysene	14.16	228	36474	4.270	ng	95
85) Indeno(1,2,3-cd)pyrene	17.27	276	15077	2.080	ng	# 94
87) Benzo(b)fluoranthene	15.22	252	39065m	5.820	ng	
88) Benzo(k)fluoranthene	15.24	252	12815m	2.005	ng	
89) Benzo(a)pyrene	15.61	252	24514	4.109	ng	# 91
91) Benzo(g,h,i)perylene	17.76	276	14844	3.003	ng	# 92

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

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