

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071318\
 Data File : BF107361.D
 Acq On : 13 Jul 2018 12:37
 Operator : JU/SJ
 Sample : J3947-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PAR-2

Manual Integrations
 APPROVED

Sohil
 7/16/2018 3:22:07 PM

Quant Time: Jul 13 14:31:36 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	55305	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	196337	20.00	ng	0.00
38) Acenaphthene-d10	9.98	164	81437	20.00	ng	0.00
63) Phenanthrene-d10	11.47	188	144839	20.00	ng	0.00
75) Chrysene-d12	14.13	240	175273	20.00	ng	0.00
86) Perylene-d12	15.67	264	161438	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	328073	100.45	ng	0.00
7) Phenol-d6	6.58	99	385611	94.54	ng	0.00
23) Nitrobenzene-d5	7.50	82	240109	67.96	ng	0.00
41) 2,4,6-Tribromophenol	10.78	330	86911	92.08	ng	0.00
44) 2-Fluorobiphenyl	9.28	172	355510	71.56	ng	0.00
78) Terphenyl-d14	13.05	244	372277	51.45	ng	0.00
Target Compounds						
10) Phenol	6.59	94	10748	2.309	ng	83
20) 3+4-Methylphenols	7.37	107	2024	Below Cal	#	81
49) Dimethylphthalate	9.68	163	66234	10.844	ng	99
70) Phenanthrene	11.49	178	18997	2.719	ng	99
74) Fluoranthene	12.69	202	33062	4.294	ng	96
77) Pyrene	12.93	202	55082	4.766	ng	98
80) Benzo(a)anthracene	14.12	228	54969	5.146	ng	# 94
82) Chrysene	14.16	228	61665	6.275	ng	# 95
85) Indeno(1,2,3-cd)pyrene	17.26	276	21086	2.528	ng	# 87
87) Benzo(b)fluoranthene	15.21	252	58012m	5.785	ng	
88) Benzo(k)fluoranthene	15.24	252	20219m	2.117	ng	
89) Benzo(a)pyrene	15.61	252	55670	6.244	ng	96
91) Benzo(a,h,i)perylene	17.75	276	20199m	2.735	ng	

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

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