

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\
 Data File : BF107469.D
 Acq On : 18 Jul 2018 11:34
 Operator : JU/SJ
 Sample : J3906-08
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 STOCKPILE-NEWSAN

Manual Integrations
 APPROVED

Sohil
 7/19/2018 2:23:17 PM

Quant Time: Jul 19 14:18:48 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 17 11:29:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	110243	20.00	ng	0.00
21) Naphthalene-d8	8.27	136	427062	20.00	ng	0.00
38) Acenaphthene-d10	10.03	164	216702	20.00	ng	0.00
63) Phenanthrene-d10	11.52	188	375620	20.00	ng	0.00
75) Chrysene-d12	14.18	240	270384m	20.00	ng	0.01
86) Perylene-d12	15.73	264	252414	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	648808	89.38	ng	0.00
7) Phenol-d6	6.61	99	893357	95.19	ng	0.00
23) Nitrobenzene-d5	7.55	82	658250	82.19	ng	0.00
41) 2,4,6-Tribromophenol	10.82	330	180922	95.02	ng	0.00
44) 2-Fluorobiphenyl	9.34	172	1014109	77.01	ng	0.00
78) Terphenyl-d14	13.11	244	763395	70.15	ng	0.00
Target Compounds						
10) Phenol	6.62	94	22253	2.212	ng	Qvalue # 73
31) Naphthalene	8.29	128	69262	3.339	ng	97
37) 2-Methylnaphthalene	8.98	142	38489	2.815	ng	# 76
49) Dimethylphthalate	9.74	163	209270	12.154	ng	# 95
51) Acenaphthene	10.06	154	47218	3.449	ng	# 70
54) Dibenzofuran	10.24	168	51211	2.421	ng	# 90
57) Fluorene	10.58	166	65438	4.670	ng	# 92
70) Phenanthrene	11.55	178	724230	38.689	ng	97
71) Anthracene	11.60	178	159222	8.540	ng	98
72) Carbazole	11.75	167	71366	4.043	ng	# 82
74) Fluoranthene	12.77	202	869906	40.886	ng	92
77) Pvrene	12.98	202	761400	41.411	ng	99
80) Benzo(a)anthracene	14.17	228	347454	21.002	ng	# 96
82) Chrysene	14.21	228	313809	19.857	ng	# 94
85) Indeno(1,2,3-cd)pyrene	17.31	276	180196	16.019	ng	# 86
87) Benzo(b)fluoranthene	15.27	252	436643m	29.554	ng	
88) Benzo(k)fluoranthene	15.29	252	120282m	8.342	ng	
89) Benzo(a)pyrene	15.67	252	294990	21.754	ng	97
90) Dibenzo(a,h)anthracene	17.32	278	47431	4.635	ng	# 88
91) Benzo(a,h,i)perylene	17.79	276	172564	16.594	ng	# 87

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

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