

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071618\
 Data File : BF107426.D
 Acq On : 17 Jul 2018 2:05
 Operator : JU/SJ
 Sample : J3975-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PAR-6

Manual Integrations
 APPROVED

Sohil
 7/17/2018 4:53:28 PM

Quant Time: Jul 17 04:27:05 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 16 17:24:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	138385	20.00	ng	0.00
21) Naphthalene-d8	8.27	136	459471	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	208242	20.00	ng	0.00
63) Phenanthrene-d10	11.52	188	394696	20.00	ng	0.00
75) Chrysene-d12	14.17	240	360737	20.00	ng	0.01
86) Perylene-d12	15.71	264	244683	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	666448	73.14	ng	0.01
7) Phenol-d6	6.61	99	829136	70.38	ng	0.01
23) Nitrobenzene-d5	7.54	82	541922	62.89	ng	0.00
41) 2,4,6-Tribromophenol	10.82	330	135612	74.12	ng	0.00
44) 2-Fluorobiphenyl	9.34	172	650177	47.24	ng	0.00
78) Terphenyl-d14	13.11	244	609469	41.98	ng	0.00
Target Compounds						
10) Phenol	6.61	94	26049	2.062	ng	# 55
31) Naphthalene	8.29	128	73191	3.279	ng	# 94
37) 2-Methylnaphthalene	8.98	142	51378	3.492	ng	# 79
49) Dimethylphthalate	9.73	163	148873	8.998	ng	# 97
51) Acenaphthene	10.06	154	67568	5.136	ng	# 94
54) Dibenzofuran	10.23	168	43736	2.151	ng	# 80
57) Fluorene	10.57	166	63777	4.737	ng	# 93
70) Phenanthrene	11.54	178	557455	28.340	ng	# 97
71) Anthracene	11.60	178	137371	7.012	ng	# 98
74) Fluoranthene	12.75	202	698046	31.223	ng	# 93
77) Pyrene	12.97	202	605439	24.681	ng	# 96
80) Benzo(a)anthracene	14.16	228	304651	13.803	ng	# 94
82) Chrysene	14.20	228	307243	14.572	ng	# 92
83) Bis(2-ethylhexyl)phthalate	14.14	149	37494	2.928	ng	# 76
85) Indeno(1,2,3-cd)pyrene	17.28	276	71170	4.742	ng	# 65
87) Benzo(b)fluoranthene	15.25	252	227966m	15.918	ng	#
88) Benzo(k)fluoranthene	15.28	252	77709m	5.560	ng	#
89) Benzo(a)pyrene	15.64	252	141035	10.729	ng	# 81
91) Benzo(a,h,i)perylene	17.75	276	68328	6.778	ng	# 82

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

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