

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070918\
 Data File : BF107229.D
 Acq On : 9 Jul 2018 15:27
 Operator : JU/SJ
 Sample : J3879-01RE
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SURCHARGE-SP-1RE

Manual Integrations
 APPROVED

Sohil
 7/10/2018 2:46:21 PM

Quant Time: Jul 09 17:28:07 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 29 14:38:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	32126	20.00	ng	-0.01
21) Naphthalene-d8	8.23	136	113096	20.00	ng	-0.02
38) Acenaphthene-d10	9.99	164	50063	20.00	ng	-0.02
63) Phenanthrene-d10	11.49	188	96190	20.00	ng	-0.01
75) Chrysene-d12	14.15	240	93053	20.00	ng	-0.02
86) Perylene-d12	15.70	264	60704	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	223879	118.00	ng	-0.01
7) Phenol-d6	6.60	99	271846	114.74	ng	-0.02
23) Nitrobenzene-d5	7.51	82	163224	80.21	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	71693	123.56	ng	-0.01
44) 2-Fluorobiphenyl	9.30	172	294403	102.59	ng	-0.02
78) Terphenyl-d14	13.07	244	346792	90.27	ng	-0.01
Target Compounds						
10) Phenol	6.61	94	5945	2.199	ng	Qvalue # 75
49) Dimethylphthalate	9.70	163	80513	21.443	ng	100
70) Phenanthrene	11.51	178	66033	14.233	ng	98
71) Anthracene	11.57	178	24158	5.101	ng	98
74) Fluoranthene	12.71	202	147073	28.761	ng	95
77) Pyrene	12.94	202	144116	23.486	ng	99
80) Benzo(a)anthracene	14.14	228	85598	15.093	ng	96
82) Chrysene	14.18	228	80056	15.344	ng	97
85) Indeno(1,2,3-cd)pyrene	17.30	276	30926	6.985	ng	# 85
87) Benzo(b)fluoranthene	15.24	252	86720m	22.997	ng	
88) Benzo(k)fluoranthene	15.27	252	26739m	7.446	ng	
89) Benzo(a)pyrene	15.64	252	51134	15.254	ng	95
90) Dibenzo(a,h)anthracene	17.30	278	9141	3.218	ng	# 69
91) Benzo(a,h,i)perylene	17.79	276	31147	11.217	ng	# 89

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

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