

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070718\
 Data File : BF107208.D
 Acq On : 7 Jul 2018 16:01
 Operator : JU/SJ
 Sample : J3879-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SURCHARGE-SP-1

Manual Integrations
 APPROVED

Sohil
 7/9/2018 2:06:29 PM

Quant Time: Jul 09 04:26:10 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	42502	20.00	ng	-0.01
21) Naphthalene-d8	8.23	136	144116	20.00	ng	-0.02
38) Acenaphthene-d10	9.99	164	62632	20.00	ng	-0.02
63) Phenanthrene-d10	11.48	188	124587	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	115568	20.00	ng	-0.02
86) Perylene-d12	15.68	264	85062	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	303324	120.85	ng	-0.01
7) Phenol-d6	6.60	99	362129	115.53	ng	-0.02
23) Nitrobenzene-d5	7.51	82	212073	81.78	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	91567	126.14	ng	-0.01
44) 2-Fluorobiphenyl	9.30	172	376015	105.11	ng	-0.02
78) Terphenyl-d14	13.07	244	461540	96.74	ng	-0.02
Target Compounds						
10) Phenol	6.61	94	8008	2.239	ng	Qvalue # 74
49) Dimethylphthalate	9.70	163	99444	21.170	ng	99
70) Phenanthrene	11.51	178	84791	14.110	ng	97
71) Anthracene	11.56	178	30947	5.046	ng	99
74) Fluoranthene	12.71	202	195332	29.492	ng	95
77) Pyrene	12.94	202	189050	24.807	ng	99
80) Benzo(a)anthracene	14.13	228	102111m	14.497	ng	
82) Chrysene	14.17	228	102741m	15.855	ng	
85) Indeno(1,2,3-cd)pyrene	17.27	276	42572	7.742	ng	# 87
87) Benzo(b)fluoranthene	15.22	252	104235m	19.727	ng	
88) Benzo(k)fluoranthene	15.25	252	35653m	7.085	ng	
89) Benzo(a)pyrene	15.62	252	69691	14.836	ng	97
90) Dibenzo(a,h)anthracene	17.27	278	11757	2.953	ng	# 76
91) Benzo(g,h,i)perylene	17.76	276	42377	10.891	ng	# 86

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

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