

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081018\
 Data File : BF108106.D
 Acq On : 11 Aug 2018 2:04
 Operator : JU/SJ
 Sample : J4409-03DL 2X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 2018-NYSEG-CLARK-AREA-10BDL

Manual Integrations
 APPROVED

Sohil
 8/13/2018 3:52:39 PM

Quant Time: Aug 11 05:45:17 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 12:24:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.02	152	259154	20.00	ng	0.00
21) Naphthalene-d8	8.31	136	906736	20.00	ng	0.00
38) Acenaphthene-d10	10.07	164	383867	20.00	ng	0.00
63) Phenanthrene-d10	11.57	188	653227	20.00	ng	0.00
75) Chrysene-d12	14.22	240	533024	20.00	ng	0.00
86) Perylene-d12	15.79	264	377636	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	843390	58.92	ng	0.01
7) Phenol-d6	6.63	99	1119434	58.85	ng	0.00
23) Nitrobenzene-d5	7.58	82	684781	42.14	ng	0.00
41) 2,4,6-Tribromophenol	10.86	330	190421	49.73	ng	0.00
44) 2-Fluorobiphenyl	9.38	172	1108263	46.35	ng	0.00
78) Terphenyl-d14	13.15	244	981299	46.81	ng	0.00
Target Compounds						
10) Phenol	6.64	94	48137	2.446	ng	Qvalue 87
31) Naphthalene	8.32	128	103675	2.514	ng	98
48) Acenaphthylene	9.93	152	88293	2.497	ng	99
49) Dimethylphthalate	9.77	163	169963	6.356	ng	# 97
51) Acenaphthene	10.10	154	109585	5.059	ng	99
57) Fluorene	10.62	166	86298	3.474	ng	96
70) Phenanthrene	11.60	178	947482	30.752	ng	99
71) Anthracene	11.64	178	363971	11.526	ng	99
74) Fluoranthene	12.79	202	1425834	42.403	ng	97
77) Pyrene	13.02	202	1511408	44.788	ng	98
80) Benzo(a)anthracene	14.21	228	579681m	18.950	ng	
82) Chrysene	14.25	228	461256	16.447	ng	97
85) Indeno(1,2,3-cd)pyrene	17.42	276	174353	7.175	ng	# 89
87) Benzo(b)fluoranthene	15.32	252	456435m	20.227	ng	
88) Benzo(k)fluoranthene	15.35	252	141770m	6.835	ng	
89) Benzo(a)pyrene	15.72	252	354881	17.563	ng	97
90) Dibenzo(a,h)anthracene	17.43	278	40729	2.436	ng	92
91) Benzo(g,h,i)perylene	17.92	276	173934	10.565	ng	# 90

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

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