

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040219\
 Data File : BF113511.D
 Acq On : 2 Apr 2019 17:50
 Operator : JU/SJ
 Sample : K2190-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VN-SP-1

Manual Integrations
 APPROVED

Sohil
 4/3/2019 12:42:28 PM

Quant Time: Apr 03 07:54:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	127111	20.00	ng	-0.01
21) Naphthalene-d8	7.97	136	471834	20.00	ng	-0.01
39) Acenaphthene-d10	9.72	164	225658	20.00	ng	-0.02
64) Phenanthrene-d10	11.20	188	406325	20.00	ng	-0.02
76) Chrysene-d12	13.83	240	295368	20.00	ng	-0.02
87) Perylene-d12	15.23	264	351044	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	710488	91.40	ng	-0.01
7) Phenol-d6	6.34	99	942133	95.80	ng	-0.02
23) Nitrobenzene-d5	7.26	82	553763	69.66	ng	-0.01
42) 2,4,6-Tribromophenol	10.51	330	262938	94.33	ng	-0.02
45) 2-Fluorobiphenyl	9.05	172	1043938	69.24	ng	-0.02
79) Terphenyl-d14	12.78	244	905131	67.54	ng	-0.02
Target Compounds						
49) Acenaphthylene	9.58	152	94429	4.025	ng	98
50) Dimethylphthalate	9.44	163	86136	5.151	ng	98
71) Phenanthrene	11.22	178	159092	7.884	ng	97
72) Anthracene	11.27	178	46545	2.271	ng	98
75) Fluoranthene	12.41	202	430665	18.880	ng	99
78) Pyrene	12.63	202	606650	29.639	ng	99
81) Benzo(a)anthracene	13.82	228	177892m	10.522	ng	
83) Chrysene	13.86	228	188313	10.965	ng	100
86) Indeno(1,2,3-cd)pyrene	16.59	276	207478	14.077	ng	95
88) Benzo(b)fluoranthene	14.84	252	318446m	14.603	ng	
89) Benzo(k)fluoranthene	14.86	252	69540m	3.631	ng	
90) Benzo(a)pyrene	15.17	252	236747	12.246	ng	99
91) Dibenzo(a,h)anthracene	16.59	278	35925m	2.149	ng	
92) Benzo(g,h,i)perylene	16.99	276	271448	16.105	ng	99

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

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