

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040219\
 Data File : BF113512.D
 Acq On : 2 Apr 2019 18:18
 Operator : JU/SJ
 Sample : K2190-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VN-SP-2

Manual Integrations
 APPROVED

Sohil
 4/3/2019 12:42:34 PM

Quant Time: Apr 03 07:55:49 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	138493	20.00	ng	-0.01
21) Naphthalene-d8	7.97	136	516439	20.00	ng	-0.01
39) Acenaphthene-d10	9.72	164	247504	20.00	ng	-0.02
64) Phenanthrene-d10	11.20	188	467154	20.00	ng	-0.02
76) Chrysene-d12	13.83	240	322774	20.00	ng	-0.01
87) Perylene-d12	15.24	264	378522	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	666594	78.70	ng	-0.01
7) Phenol-d6	6.34	99	894829	83.51	ng	-0.02
23) Nitrobenzene-d5	7.26	82	529905	60.90	ng	-0.01
42) 2,4,6-Tribromophenol	10.51	330	263330	86.13	ng	-0.02
45) 2-Fluorobiphenyl	9.05	172	1001797	57.80	ng	-0.02
79) Terphenyl-d14	12.78	244	903088	61.67	ng	-0.02
Target Compounds						
49) Acenaphthylene	9.58	152	219711	8.538	ng	98
50) Dimethylphthalate	9.44	163	64170	3.499	ng	99
71) Phenanthrene	11.22	178	285175	12.293	ng	99
72) Anthracene	11.27	178	60262	2.557	ng	98
75) Fluoranthene	12.41	202	855005	32.601	ng	99
78) Pyrene	12.64	202	1229215	54.956	ng	100
81) Benzo(a)anthracene	13.82	228	423213m	22.906	ng	
83) Chrysene	13.86	228	388160	20.683	ng	99
86) Indeno(1,2,3-cd)pyrene	16.59	276	373385	23.183	ng	96
88) Benzo(b)fluoranthene	14.84	252	578298m	24.594	ng	
89) Benzo(k)fluoranthene	14.86	252	159012m	7.701	ng	
90) Benzo(a)pyrene	15.18	252	507052	24.324	ng	99
91) Dibenzo(a,h)anthracene	16.59	278	68340m	3.792	ng	
92) Benzo(g,h,i)perylene	17.00	276	451939	24.867	ng	99

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

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