

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071819\
 Data File : BF115661.D
 Acq On : 18 Jul 2019 23:12
 Operator : HP/JU
 Sample : K3875-02 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HJDC-S2

Manual Integrations
 APPROVED

mohammad
 7/19/2019 3:58:52 PM

Quant Time: Jul 19 08:14:31 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 14:03:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	150436	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	598452	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	260397	20.00	ng	-0.01
64) Phenanthrene-d10	11.40	188	412174	20.00	ng	0.00
76) Chrysene-d12	14.04	240	344509	20.00	ng	-0.01
87) Perylene-d12	15.51	264	280379	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	346363	37.66	ng	0.00
7) Phenol-d6	6.50	99	482036	41.31	ng	-0.01
23) Nitrobenzene-d5	7.44	82	305375	29.67	ng	-0.01
42) 2,4,6-Tribromophenol	10.70	330	100869	33.19	ng	-0.01
45) 2-Fluorobiphenyl	9.23	172	483747	32.52	ng	-0.02
79) Terphenyl-d14	12.98	244	424969	22.76	ng	-0.02

Target Compounds				Qvalue		
31) Naphthalene	8.19	128	114802	3.961	ng	96
50) Dimethylphthalate	9.62	163	66445	3.392	ng	99
52) Acenaphthene	9.95	154	67519	4.164	ng	98
55) Dibenzofuran	10.12	168	70867	3.077	ng	# 96
58) Fluorene	10.46	166	102265	6.212	ng	99
71) Phenanthrene	11.43	178	698183	33.046	ng	96
72) Anthracene	11.47	178	187289	8.578	ng	98
73) Carbazole	11.63	167	91289	4.768	ng	98
75) Fluoranthene	12.62	202	955649	42.804	ng	98
78) Pyrene	12.84	202	823410	28.210	ng	97
81) Benzo(a)anthracene	14.03	228	487737	21.490	ng	99
83) Chrysene	14.06	228	414122	18.176	ng	96
86) Indeno(1,2,3-cd)pyrene	16.98	276	162493	8.395	ng	95
88) Benzo(b)fluoranthene	15.08	252	424164m	23.494	ng	
89) Benzo(k)fluoranthene	15.10	252	108782m	6.758	ng	
90) Benzo(a)pyrene	15.44	252	259772	16.741	ng	97
91) Dibenzo(a,h)anthracene	16.99	278	45870	3.367	ng	# 87
92) Benzo(g,h,i)perylene	17.42	276	140242	10.322	ng	97

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

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