

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\
 Data File : BF117612.D
 Acq On : 30 Oct 2019 1:42
 Operator : JU
 Sample : K5576-19 5X
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CSA-2-1-8

Manual Integrations
 APPROVED

mohammad
 10/30/2019 4:53:43 PM

Quant Time: Oct 30 07:06:04 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 18:53:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	39139	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	161238	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	73978	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	108502	20.00	ng	0.00
76) Chrysene-d12	13.90	240	70359	20.00	ng	0.00
87) Perylene-d12	15.33	264	57764	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.37	112	32051	12.18	ng	0.02
7) Phenol-d6	6.40	99	52328	14.81	ng	0.02
23) Nitrobenzene-d5	7.30	82	28482	8.72	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	7899	12.34	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	53578	10.20	ng	-0.01
79) Terphenyl-d14	12.83	244	43368	10.05	ng	0.00

Target Compounds

						Qvalue
31) Naphthalene	8.03	128	39078	4.890	ng	97
37) 2-Methylnaphthalene	8.73	142	18915	3.512	ng	99
38) 1-Methylnaphthalene	8.82	142	19783	3.909	ng	100
49) Acenaphthylene	9.63	152	13566	2.000	ng	97
52) Acenaphthene	9.79	154	53774	12.922	ng	97
55) Dibenzofuran	9.97	168	46112	7.648	ng	99
58) Fluorene	10.31	166	53904	10.812	ng	99
71) Phenanthrene	11.28	178	677845	107.589	ng	99
72) Anthracene	11.33	178	178459	27.606	ng	99
73) Carbazole	11.49	167	41438	7.014	ng	97
75) Fluoranthene	12.47	202	585165	92.779	ng	99
78) Pvrene	12.70	202	543711	89.967	ng	100
80) Butylbenzylphthalate	13.30	149	9564	3.180	ng	# 92
81) Benzo(a)anthracene	13.89	228	189712	39.973	ng	98
83) Chrysene	13.93	228	162097	34.849	ng	98
86) Indeno(1,2,3-cd)pyrene	16.75	276	48916	14.086	ng	96
88) Benzo(b)fluoranthene	14.93	252	135510m	39.727	ng	
89) Benzo(k)fluoranthene	14.94	252	51029m	14.596	ng	
90) Benzo(a)pvrene	15.27	252	99073	31.945	ng	95
91) Dibenzo(a,h)anthracene	16.75	278	13583m	5.065	ng	
92) Benzo(g,h,i)perylene	17.18	276	46955	18.506	ng	97

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

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