

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF093019\
 Data File : BF117270.D
 Acq On : 30 Sep 2019 19:35
 Operator : JU
 Sample : K5093-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-01-COMP

Manual Integrations
 APPROVED

mohammad
 10/2/2019 8:32:41 AM

Quant Time: Oct 01 09:38:49 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 13 16:23:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	27845	20.00	ng	-0.01
21) Naphthalene-d8	8.09	136	114205	20.00	ng	-0.01
39) Acenaphthene-d10	9.84	164	57908	20.00	ng	-0.02
64) Phenanthrene-d10	11.33	188	94628	20.00	ng	0.00
76) Chrysene-d12	13.97	240	64876	20.00	ng	0.00
87) Perylene-d12	15.43	264	72718	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	159075	89.19	ng	0.00
7) Phenol-d6	6.45	99	223691	97.05	ng	0.00
23) Nitrobenzene-d5	7.37	82	139623	64.24	ng	-0.02
42) 2,4,6-Tribromophenol	10.63	330	39792	82.22	ng	-0.01
45) 2-Fluorobiphenyl	9.16	172	253293	68.39	ng	-0.01
79) Terphenyl-d14	12.91	244	235002	67.71	ng	0.00
Target Compounds						
10) Phenol	6.46	94	6995	2.753	ng	Qvalue # 79
31) Naphthalene	8.11	128	48649	8.857	ng	98
37) 2-Methylnaphthalene	8.81	142	8866	2.397	ng	99
38) 1-Methylnaphthalene	8.90	142	8199	2.367	ng	# 100
50) Dimethylphthalate	9.56	163	34099	8.316	ng	99
52) Acenaphthene	9.87	154	30445	9.564	ng	99
55) Dibenzofuran	10.05	168	21408	4.558	ng	96
58) Fluorene	10.39	166	21985	5.811	ng	# 98
71) Phenanthrene	11.35	178	306017	56.935	ng	98
72) Anthracene	11.40	178	86031	15.678	ng	97
73) Carbazole	11.56	167	26530	5.094	ng	97
75) Fluoranthene	12.55	202	343492	62.197	ng	99
78) Pvrene	12.77	202	301103	55.313	ng	98
81) Benzo(a)anthracene	13.96	228	110576	25.511	ng	98
83) Chrysene	13.99	228	94057	22.249	ng	96
86) Indeno(1,2,3-cd)pyrene	16.87	276	56702	18.385	ng	95
88) Benzo(b)fluoranthene	15.00	252	117030m	26.569	ng	
89) Benzo(k)fluoranthene	15.03	252	39063m	9.362	ng	
90) Benzo(a)pvrene	15.36	252	93193	24.110	ng	98
91) Dibenzo(a,h)anthracene	16.87	278	13267m	4.309	ng	
92) Benzo(g,h,i)perylene	17.31	276	58436	19.044	ng	98

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

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