

Data Path : Z:\HPCHEM1\BNA F\DATA\BF033117\
 Data File : BF094110.D
 Acq On : 31 Mar 2017 21:43
 Operator : SJ/MA
 Sample : I2463-10
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-BB-7-8-SSW

Manual Integrations
 APPROVED

mohammad
 4/3/2017 3:26:27 PM

Quant Time: Apr 01 00:43:54 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 31 14:38:29 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.90	152	186945	20.00	ng	0.00
21) Naphthalene-d8	7.41	136	719740	20.00	ng	0.00
38) Acenaphthene-d10	9.55	164	307930	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	453952	20.00	ng	0.00
75) Chrysene-d12	14.63	240	305220	20.00	ng	0.00
86) Perylene-d12	16.26	264	300846	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	4.46	112	1464322	132.30	ng	0.02
7) Phenol-d6	5.53	99	1875377	136.96	ng	0.02
23) Nitrobenzene-d5	6.56	82	1062039	84.21	ng	0.00
41) 2,4,6-Tribromophenol	10.53	330	263141	108.14	ng	0.00
44) 2-Fluorobiphenyl	8.73	172	1623506	80.42	ng	0.00
78) Terphenyl-d14	13.36	244	835192	61.34	ng	0.00
Target Compounds						
9) Benzaldehyde	5.40	77	20470	2.21	ng	Qvalue # 81
10) Phenol	5.54	94	75681	4.86	ng	88
17) 2-Methylphenol	6.22	107	99823	9.58	ng	99
20) 3+4-Methylphenols	6.40	107	210960	16.72	ng	# 60
27) 2,4-Dimethylphenol	7.02	122	296655	29.02	ng	98
31) Naphthalene	7.43	128	1286171	37.16	ng	99
37) 2-Methylnaphthalene	8.27	142	371732	16.11	ng	98
48) Acenaphthylene	9.37	152	114430	3.54	ng	# 95
49) Dimethylphthalate	9.23	163	117392	5.36	ng	99
51) Acenaphthene	9.59	154	424981	22.54	ng	99
54) Dibenzofuran	9.80	168	502280	19.57	ng	99
57) Fluorene	10.22	166	687183	32.29	ng	98
70) Phenanthrene	11.42	178	3184788	126.12	ng	100
71) Anthracene	11.47	178	895236	34.43	ng	99
72) Carbazole	11.67	167	803130	30.12	ng	98
74) Fluoranthene	12.88	202	2165009	83.73	ng	100
77) Pyrene	13.16	202	2083037	94.04	ng	99
80) Benzo(a)anthracene	14.62	228	923379	51.88	ng	99
82) Chrysene	14.67	228	913555	55.31	ng	96
85) Indeno(1,2,3-cd)pyrene	17.62	276	363072	25.38	ng	93
87) Benzo(b)fluoranthene	15.86	252	1040846m	56.44	ng	
88) Benzo(k)fluoranthene	15.89	252	247071m	13.69	ng	
89) Benzo(a)pyrene	16.21	252	743121	43.76	ng	97
90) Dibenzo(a,h)anthracene	17.63	278	89926m	6.25	ng	
91) Benzo(g,h,i)perylene	18.02	276	328747	22.68	ng	98

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

