

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091317\
 Data File : BF098505.D
 Acq On : 13 Sep 2017 22:36
 Operator : SJ/JU
 Sample : I5160-02
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G-6-1

Manual Integrations
 APPROVED

Sohil
 9/14/2017 5:52:52 PM

Quant Time: Sep 14 07:25:52 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF091117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 12 02:14:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	184812	20.00	ng	-0.01
21) Naphthalene-d8	7.94	136	766200	20.00	ng	-0.01
38) Acenaphthene-d10	9.69	164	344566	20.00	ng	-0.02
63) Phenanthrene-d10	11.17	188	554652	20.00	ng	-0.02
75) Chrysene-d12	13.81	240	356700	20.00	ng	0.00
86) Perylene-d12	15.21	264	327123	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	1258143	100.65	ng	0.00
7) Phenol-d6	6.32	99	1589052	100.13	ng	0.00
23) Nitrobenzene-d5	7.23	82	917643	66.77	ng	-0.01
41) 2,4,6-Tribromophenol	10.48	330	248010	107.84	ng	-0.01
44) 2-Fluorobiphenyl	9.02	172	1563646	76.92	ng	-0.02
78) Terphenyl-d14	12.76	244	982557	59.46	ng	-0.01

Target Compounds				Qvalue		
10) Phenol	6.33	94	54588	2.961	ng	# 67
31) Naphthalene	7.96	128	333876	8.815	ng	99
37) 2-Methylnaphthalene	8.65	142	236841	10.323	ng	99
48) Acenaphthylene	9.55	152	179174	5.471	ng	99
49) Dimethylphthalate	9.41	163	347393	13.025	ng	99
54) Dibenzofuran	9.89	168	115821	4.223	ng	# 97
70) Phenanthrene	11.19	178	406288	13.818	ng	99
71) Anthracene	11.25	178	186102	6.165	ng	97
72) Carbazole	11.40	167	66031	2.347	ng	99
74) Fluoranthene	12.38	202	739037	25.107	ng	99
77) Pyrene	12.61	202	703972	25.018	ng	98
79) Butylbenzylphthalate	13.23	149	30022	2.459	ng	# 76
80) Benzo(a)anthracene	13.80	228	423669m	20.259	ng	
82) Chrysene	13.83	228	499686	23.638	ng	97
83) Bis(2-ethylhexyl)phthalate	13.79	149	121441	7.927	ng	99
85) Indeno(1,2,3-cd)pyrene	16.55	276	228184	15.386	ng	# 91
87) Benzo(b)fluoranthene	14.82	252	766152m	37.249	ng	
88) Benzo(k)fluoranthene	14.84	252	242361m	12.622	ng	
89) Benzo(a)pyrene	15.15	252	347240	18.950	ng	96
90) Dibenzo(a,h)anthracene	16.55	278	62167	4.664	ng	# 76
91) Benzo(g,h,i)perylene	16.96	276	191368	14.273	ng	96

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

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