

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110517\
 Data File : BF100217.D
 Acq On : 5 Nov 2017 12:16
 Operator : SJ/JU
 Sample : I6299-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-02-110217

Manual Integrations
 APPROVED

Sohil
 11/6/2017 2:33:19 PM

Quant Time: Nov 06 06:13:46 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 03 15:46:10 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	93933	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	376873	20.00	ng	0.00
38) Acenaphthene-d10	9.80	164	164946	20.00	ng	-0.01
63) Phenanthrene-d10	11.29	188	258908	20.00	ng	-0.01
75) Chrysene-d12	13.94	240	165092	20.00	ng	-0.01
86) Perylene-d12	15.40	264	184042	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.39	112	570682	95.38	ng	0.00
7) Phenol-d6	6.42	99	666056	93.89	ng	-0.01
23) Nitrobenzene-d5	7.33	82	367997	62.86	ng	-0.02
41) 2,4,6-Tribromophenol	10.59	330	131902	87.75	ng	-0.02
44) 2-Fluorobiphenyl	9.12	172	722798	67.61	ng	-0.01
78) Terphenyl-d14	12.87	244	464512	61.49	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
49) Dimethylphthalate	9.52	163	95900	7.342	ng	99
51) Acenaphthene	9.83	154	27962	2.742	ng	95
57) Fluorene	10.35	166	56874	4.772	ng	96
70) Phenanthrene	11.32	178	529732	36.255	ng	98
71) Anthracene	11.37	178	129708	8.745	ng	97
72) Carbazole	11.54	167	47173	3.382	ng	97
74) Fluoranthene	12.51	202	587043	38.660	ng	97
77) Pyrene	12.74	202	523921	41.735	ng	99
80) Benzo(a)anthracene	13.93	228	208799	19.951	ng	96
82) Chrysene	13.97	228	211072	21.452	ng	98
85) Indeno(1,2,3-cd)pyrene	16.89	276	64313	7.120	ng	95
87) Benzo(b)fluoranthene	14.98	252	196915m	16.944	ng	
88) Benzo(k)fluoranthene	15.01	252	73819m	6.736	ng	
89) Benzo(a)pyrene	15.34	252	136106	13.038	ng	99
90) Dibenzo(a,h)anthracene	16.87	278	20383	2.197	ng	# 82
91) Benzo(g,h,i)perylene	17.33	276	59532	6.560	ng	93

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

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