

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091015\
 Data File : BF081574.D
 Acq On : 10 Sep 2015 19:06
 Operator : UM/IZ
 Sample : G3593-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP(E2PM-MLSP1-SP12)

Quant Time: Sep 11 01:33:49 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 11 00:38:02 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.26	152	49135	20.00	ng	0.00
21) Naphthalene-d8	11.16	136	186206	20.00	ng	-0.01
38) Acenaphthene-d10	15.29	164	79736	20.00	ng	-0.01
63) Phenanthrene-d10	18.81	188	129559	20.00	ng	0.00
75) Chrysene-d12	23.42	240	104266	20.00	ng	0.00
86) Perylene-d12	24.86	264	89021	20.00	ng	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.43	112	342613	108.19	ng	0.01
7) Phenol-d6	7.69	99	464790	110.88	ng	0.00
23) Nitrobenzene-d5	9.57	82	308180	79.85	ng	0.00
41) 2,4,6-Tribromophenol	17.20	330	71088	100.13	ng	-0.01
44) 2-Fluorobiphenyl	13.81	172	432825	69.56	ng	-0.01
78) Terphenyl-d14	22.15	244	244298	54.54	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
37) 2-Methylnaphthalene	12.86	142	19126	2.96	ng	# 88
48) Acenaphthylene	14.95	152	49432	5.26	ng	97
49) Dimethylphthalate	14.87	163	118027	19.05	ng	# 97
51) Acenaphthene	15.37	154	27189	5.09	ng	# 79
57) Fluorene	16.61	166	54521	9.20	ng	# 95
70) Phenanthrene	18.87	178	547983	76.08	ng	97
71) Anthracene	18.98	178	133684	18.07	ng	97
72) Carbazole	19.46	167	27780	4.00	ng	# 90
73) Di-n-butylphthalate	20.52	149	87523	10.67	ng	# 98
74) Fluoranthene	21.49	202	712878	91.43	ng	85
77) Pyrene	21.83	202	770518	99.76	ng	# 92
80) Benzo(a)anthracene	23.41	228	357124	53.78	ng	# 92
82) Chrysene	23.45	228	332479	55.86	ng	92
83) Bis(2-ethylhexyl)phthalate	23.52	149	11873	2.68	ng	# 92
85) Indeno(1,2,3-cd)pyrene	25.93	276	117851	26.99	ng	# 79
87) Benzo(b)fluoranthene	24.53	252	323223m	50.86	ng	
88) Benzo(k)fluoranthene	24.53	252	149294m	28.31	ng	
89) Benzo(a)pyrene	24.81	252	270265	50.18	ng	# 84
90) Dibenzo(a,h)anthracene	25.93	278	38260	9.19	ng	# 79
91) Benzo(g,h,i)perylene	26.24	276	112935	28.43	ng	# 73

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

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