

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF011218\
 Data File : BF102059.D
 Acq On : 12 Jan 2018 22:39
 Operator : SJ/JU
 Sample : J1015-03
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-011118-B

Manual Integrations
 APPROVED

Sohil
 1/15/2018 4:22:23 PM

Quant Time: Jan 13 01:07:54 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 11 12:53:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	233362	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	874855	20.00	ng	0.00
38) Acenaphthene-d10	9.76	164	325159	20.00	ng	0.00
63) Phenanthrene-d10	11.25	188	472801	20.00	ng	0.00
75) Chrysene-d12	13.89	240	384154	20.00	ng	0.00
86) Perylene-d12	15.32	264	302834	20.00	ng	0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.34	112	1478255	89.45	ng	0.01
7) Phenol-d6	6.37	99	1767938	89.66	ng	0.00
23) Nitrobenzene-d5	7.30	82	1066148	71.61	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	256312	90.33	ng	0.00
44) 2-Fluorobiphenyl	9.09	172	1619085	77.79	ng	0.00
78) Terphenyl-d14	12.84	244	1102917	59.61	ng	0.00

Target Compounds

						Qvalue
10) Phenol	6.38	94	88322	3.964	ng	98
31) Naphthalene	8.03	128	144354	3.302	ng	99
37) 2-Methylnaphthalene	8.72	142	113840	3.956	ng	99
48) Acenaphthylene	9.63	152	329379	9.313	ng	100
49) Dimethylphthalate	9.48	163	197633	7.577	ng	99
57) Fluorene	10.31	166	118697	5.828	ng	97
70) Phenanthrene	11.27	178	539321	19.528	ng	99
71) Anthracene	11.32	178	250877	8.957	ng	99
74) Fluoranthene	12.46	202	655464	24.123	ng	98
77) Pyrene	12.69	202	813223	23.949	ng	99
80) Benzo(a)anthracene	13.87	228	432353m	17.640	ng	
82) Chrysene	13.91	228	438472	17.156	ng	98
85) Indeno(1,2,3-cd)pyrene	16.70	276	166149	8.932	ng	99
87) Benzo(b)fluoranthene	14.91	252	423215m	21.433	ng	
88) Benzo(k)fluoranthene	14.93	252	160316m	8.416	ng	
89) Benzo(a)pyrene	15.26	252	317144	17.849	ng	# 96
90) Dibenzo(a,h)anthracene	16.71	278	46690	3.293	ng	# 76
91) Benzo(g,h,i)perylene	17.12	276	176471	12.355	ng	95

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

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