

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102341.D
 Acq On : 23 Jan 2018 13:11
 Operator : SJ/JU
 Sample : J1239-13
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-11R-01

Manual Integrations
 APPROVED

Sohil
 1/24/2018 2:46:42 PM

Quant Time: Jan 24 11:31:47 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 23 00:43:41 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	160469	20.00	ng	0.01
21) Naphthalene-d8	8.02	136	473079	20.00	ng	0.02
38) Acenaphthene-d10	9.76	164	291282	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	493381	20.00	ng	0.00
75) Chrysene-d12	13.87	240	400819	20.00	ng	0.00
86) Perylene-d12	15.30	264	300480	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1189116	112.36	ng	0.02
7) Phenol-d6	6.43	99	768195m	60.21	ng	0.07
23) Nitrobenzene-d5	7.32	82	977099	118.69	ng	0.04
41) 2,4,6-Tribromophenol	10.55	330	319676	110.76	ng	0.00
44) 2-Fluorobiphenyl	9.09	172	1579062	79.71	ng	0.00
78) Terphenyl-d14	12.83	244	1555750	87.50	ng	0.00
Target Compounds						Qvalue
3) Pyridine	3.19	79	306877	23.353	ng	96
10) Phenol	6.49	94	16258610m	1167.220	ng	
17) 2-Methylphenol	7.04	107	3224072	334.621	ng	97
20) 3+4-Methylphenols	7.23	107	9789455	863.893	ng	88
27) 2,4-Dimethylphenol	7.70	122	1394210	216.548	ng	97
31) Naphthalene	8.05	128	10301355	436.127	ng	97
37) 2-Methylnaphthalene	8.73	142	593328	37.719	ng	100
45) 1,1'-Biphenyl	9.18	154	141779	5.444	ng	98
48) Acenaphthylene	9.62	152	1392869	44.162	ng	99
54) Dibenzofuran	9.96	168	405991	16.286	ng	99
57) Fluorene	10.30	166	315139	15.953	ng	# 96
70) Phenanthrene	11.26	178	425298	14.793	ng	98
71) Anthracene	11.31	178	131943m	4.496	ng	
72) Carbazole	11.49	167	1580505	55.207	ng	99
76) Benzidine	12.58	184	65202	4.842	ng	97

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

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