

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012218\
 Data File : BF102326.D
 Acq On : 23 Jan 2018 5:37
 Operator : SJ/JU
 Sample : J1214-02DL 2X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-PH3-MGP-02DL

Manual Integrations
 APPROVED

Sohil
 1/23/2018 12:47:06 PM

Quant Time: Jan 23 08:04:35 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	6.72	152	128171	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	499285	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	203257	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	293433	20.00	ng	0.00
75) Chrysene-d12	13.88	240	241681	20.00	ng	0.00
86) Perylene-d12	15.31	264	211509	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	101104	11.96	ng	0.00
7) Phenol-d6	6.36	99	125530	12.32	ng	0.00
23) Nitrobenzene-d5	7.27	82	65740	7.57	ng	-0.01
41) 2,4,6-Tribromophenol	10.55	330	18016	8.95	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	80585	5.83	ng	0.00
78) Terphenyl-d14	12.82	244	53400	4.98	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.02	128	827319	33.188	ng	99
37) 2-Methylnaphthalene	8.71	142	175118	10.548	ng	99
45) 1,1'-Biphenyl	9.17	154	122153	6.721	ng	97
48) Acenaphthylene	9.62	152	505017	22.946	ng	99
51) Acenaphthene	9.79	154	111945	8.709	ng	99
54) Dibenzofuran	9.96	168	53881	3.097	ng	# 84
57) Fluorene	10.30	166	369565	26.810	ng	99
70) Phenanthrene	11.27	178	1691686	98.935	ng	99
71) Anthracene	11.32	178	496036	28.422	ng	99
74) Fluoranthene	12.46	202	877912	50.348	ng	99
77) Pyrene	12.69	202	1146858	61.376	ng	100
80) Benzo(a)anthracene	13.87	228	326484m	21.942	ng	
82) Chrysene	13.90	228	301204	19.934	ng	99
85) Indeno(1,2,3-cd)pyrene	16.69	276	88629	7.102	ng	94
87) Benzo(b)fluoranthene	14.90	252	247129m	18.027	ng	
88) Benzo(k)fluoranthene	14.92	252	77253m	5.965	ng	
89) Benzo(a)pyrene	15.24	252	235263	19.028	ng	98
90) Dibenz(a,h)anthracene	16.70	278	21518	2.148	ng	94
91) Benzo(a,h,i)perylene	17.12	276	98893	10.000	ng	100

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

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