

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030818\
 Data File : BF103522.D
 Acq On : 8 Mar 2018 17:48
 Operator : SJ/JU
 Sample : J1819-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-9

Manual Integrations
 APPROVED

Sohil
 3/9/2018 11:39:28 AM

Quant Time: Mar 09 00:28:32 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 08 14:20:19 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	141999	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	579353	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	231462	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	317712	20.00	ng	0.00
75) Chrysene-d12	14.07	240	208229	20.00	ng	0.00
86) Perylene-d12	15.58	264	191003	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	783554	83.46	ng	0.01
7) Phenol-d6	6.52	99	1054740	91.81	ng	0.01
23) Nitrobenzene-d5	7.43	82	648997	63.97	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	142365	72.66	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	891642	62.16	ng	0.00
78) Terphenyl-d14	12.99	244	518262	59.83	ng	0.00
Target Compounds						Qvalue
51) Acenaphthene	9.94	154	189801	13.787	ng	98
54) Dibenzofuran	10.12	168	73998	3.916	ng	98
57) Fluorene	10.46	166	43191	2.683	ng	# 94
70) Phenanthrene	11.43	178	244688	13.994	ng	99
71) Anthracene	11.48	178	224987	12.548	ng	98
73) Di-n-butylphthalate	11.95	149	240867	10.781	ng	99
74) Fluoranthene	12.64	202	821303	46.744	ng	99
77) Pyrene	12.86	202	767591	47.281	ng	99
80) Benzo(a)anthracene	14.06	228	320848m	23.748	ng	
82) Chrysene	14.09	228	320234m	24.411	ng	
83) Bis(2-ethylhexyl)phthalate	14.01	149	257350	22.233	ng	100
85) Indeno(1,2,3-cd)pyrene	17.12	276	140763	13.554	ng	92
87) Benzo(b)fluoranthene	15.13	252	350560m	27.915	ng	
88) Benzo(k)fluoranthene	15.16	252	112569m	9.519	ng	
89) Benzo(a)pyrene	15.52	252	241659	21.549	ng	# 90
90) Dibenzo(a,h)anthracene	17.12	278	33714	3.891	ng	# 68
91) Benzo(a,h,i)perylene	17.60	276	130614	15.422	ng	91

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

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