

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030118\
 Data File : BF103329.D
 Acq On : 1 Mar 2018 11:55
 Operator : SJ/JU
 Sample : J1709-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DRUM-1-2

Manual Integrations
 APPROVED

Sohil
 3/2/2018 3:30:20 PM

Quant Time: Mar 01 18:19:14 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 17:05:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	161997	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	655606	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	287102	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	464019	20.00	ng	0.00
75) Chrysene-d12	14.06	240	219976	20.00	ng	0.00
86) Perylene-d12	15.57	264	270366	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.50	112	1037528	96.87	ng	0.01
7) Phenol-d6	6.51	99	1310272	99.97	ng	0.00
23) Nitrobenzene-d5	7.44	82	799128	69.61	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	194860	80.18	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1027433	56.82	ng	0.00
78) Terphenyl-d14	12.99	244	647978	70.81	ng	0.00

Target Compounds

						Qvalue
31) Naphthalene	8.17	128	176370	5.495	ng	99
37) 2-Methylnaphthalene	8.87	142	71519	3.448	ng	96
49) Dimethylphthalate	9.62	163	118713	5.154	ng	99
51) Acenaphthene	9.95	154	134975	7.904	ng	98
54) Dibenzofuran	10.12	168	148355	6.329	ng	92
57) Fluorene	10.46	166	204586	10.245	ng	97
70) Phenanthrene	11.44	178	1494995	58.544	ng	99
71) Anthracene	11.48	178	488566	18.658	ng	99
72) Carbazole	11.64	167	154997	5.944	ng	98
74) Fluoranthene	12.63	202	1528190	59.552	ng	99
77) Pyrene	12.86	202	1371614	79.974	ng	100
80) Benzo(a)anthracene	14.05	228	643654	45.096	ng	98
82) Chrysene	14.09	228	561169	40.493	ng	97
83) Bis(2-ethylhexyl)phthalate	14.02	149	43877	3.588	ng	# 97
85) Indeno(1,2,3-cd)pyrene	17.10	276	321108	29.267	ng	96
87) Benzo(b)fluoranthene	15.13	252	755978m	42.527	ng	
88) Benzo(k)fluoranthene	15.16	252	210975m	12.604	ng	
89) Benzo(a)pyrene	15.51	252	560576	35.313	ng	# 95
90) Dibenzo(a,h)anthracene	17.10	278	89314m	7.281	ng	
91) Benzo(a,h,i)perylene	17.56	276	304190	25.374	ng	96

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

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