

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030518\
 Data File : BF103407.D
 Acq On : 5 Mar 2018 14:10
 Operator : SJ/JU
 Sample : J1751-03 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-030118-B

Manual Integrations
 APPROVED

Sohil
 3/6/2018 4:22:51 PM

Quant Time: Mar 05 15:25:42 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 05 13:40:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	138143	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	571680	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	243629	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	365997	20.00	ng	0.00
75) Chrysene-d12	14.07	240	228833	20.00	ng	0.00
86) Perylene-d12	15.59	264	196131	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	535398	58.62	ng	0.00
7) Phenol-d6	6.52	99	641082	57.36	ng	0.00
23) Nitrobenzene-d5	7.43	82	367525	36.71	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	86264	41.83	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	544934	32.55	ng	0.00
78) Terphenyl-d14	12.99	244	359275	37.74	ng	0.00
Target Compounds						
10) Phenol	6.53	94	27049	2.085	ng	Qvalue # 54
48) Acenaphthylene	9.77	152	80924	3.257	ng	98
49) Dimethylphthalate	9.62	163	127523	6.525	ng	98
51) Acenaphthene	9.95	154	75463	5.208	ng	99
54) Dibenzofuran	10.12	168	61884	3.111	ng	98
57) Fluorene	10.46	166	110959	6.548	ng	98
70) Phenanthrene	11.44	178	853049	42.352	ng	99
71) Anthracene	11.49	178	343625	16.637	ng	99
74) Fluoranthene	12.64	202	1489562	73.593	ng	99
77) Pyrene	12.87	202	1371474	76.871	ng	99
80) Benzo(a)anthracene	14.06	228	743925	50.104	ng	97
82) Chrysene	14.10	228	602128	41.766	ng	98
85) Indeno(1,2,3-cd)pyrene	17.14	276	257442	22.556	ng	95
87) Benzo(b)fluoranthene	15.14	252	785938m	60.947	ng	
88) Benzo(k)fluoranthene	15.17	252	237427m	19.552	ng	
89) Benzo(a)pyrene	15.53	252	522581	45.380	ng	95
90) Dibenzo(a,h)anthracene	17.13	278	71079	7.988	ng	# 82
91) Benzo(g,h,i)perylene	17.61	276	228550	26.280	ng	93

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

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