

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072318\
 Data File : BF107592.D
 Acq On : 23 Jul 2018 17:05
 Operator : JU/SJ
 Sample : J4030-11 2X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB8-12-0

Manual Integrations
 APPROVED

Sohil
 7/24/2018 5:26:12 PM

Quant Time: Jul 23 19:05:21 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 20 16:29:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	317882	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	1149516	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	435628	20.00	ng	0.00
63) Phenanthrene-d10	11.52	188	792289	20.00	ng	0.00
75) Chrysene-d12	14.17	240	566122	20.00	ng	0.00
86) Perylene-d12	15.72	264	438084	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	861600	43.58	ng	0.00
7) Phenol-d6	6.61	99	1094901	46.12	ng	0.00
23) Nitrobenzene-d5	7.54	82	631735	37.09	ng	-0.01
41) 2,4,6-Tribromophenol	10.81	330	217915	43.99	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	1100321	35.81	ng	0.00
78) Terphenyl-d14	13.10	244	945647	42.76	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.72	163	239491	7.294	ng	98
70) Phenanthrene	11.54	178	700933	18.155	ng	99
71) Anthracene	11.59	178	175060	4.503	ng	98
72) Carbazole	11.75	167	87118	2.155	ng	99
74) Fluoranthene	12.74	202	1413673	34.122	ng	97
77) Pyrene	12.97	202	1258836	32.508	ng	98
80) Benzo(a)anthracene	14.16	228	535450	16.140	ng	97
82) Chrysene	14.19	228	527902	16.565	ng	97
85) Indeno(1,2,3-cd)pyrene	17.31	276	259739	9.617	ng	# 89
87) Benzo(b)fluoranthene	15.25	252	570046m	23.776	ng	
88) Benzo(k)fluoranthene	15.28	252	232748m	9.909	ng	
89) Benzo(a)pyrene	15.65	252	387248	17.657	ng	93
90) Dibenzo(a,h)anthracene	17.31	278	63638	3.357	ng	# 69
91) Benzo(a,h,i)perylene	17.79	276	266368	14.244	ng	91

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

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