

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF072618\
 Data File : BF107720.D
 Acq On : 27 Jul 2018 00:31
 Operator : JU/SJ
 Sample : J4121-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-9(3-5)

Manual Integrations
 APPROVED

Sohil
 7/27/2018 2:54:17 PM

Quant Time: Jul 27 06:00:08 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	6.96	152	206815	20.00	ng	-0.02
21) Naphthalene-d8	8.24	136	853139	20.00	ng	-0.02
38) Acenaphthene-d10	10.00	164	361455	20.00	ng	-0.02
63) Phenanthrene-d10	11.49	188	567395	20.00	ng	-0.02
75) Chrysene-d12	14.15	240	447013	20.00	ng	-0.02
86) Perylene-d12	15.68	264	508057	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	1278140	99.36	ng	0.00
7) Phenol-d6	6.60	99	1535683	99.43	ng	-0.01
23) Nitrobenzene-d5	7.52	82	1010877	79.97	ng	-0.02
41) 2,4,6-Tribromophenol	10.80	330	389997	94.88	ng	-0.02
44) 2-Fluorobiphenyl	9.31	172	1400791	61.28	ng	-0.02
78) Terphenyl-d14	13.08	244	914219	52.36	ng	-0.02
Target Compounds						Qvalue
31) Naphthalene	8.27	128	231503	6.171	ng	98
37) 2-Methylnaphthalene	8.95	142	72757	2.799	ng	97
48) Acenaphthylene	9.86	152	113783	3.147	ng	98
49) Dimethylphthalate	9.71	163	127461	4.679	ng	98
51) Acenaphthene	10.03	154	138751	6.125	ng	96
54) Dibenzofuran	10.21	168	161993	4.857	ng	99
57) Fluorene	10.55	166	262650	11.039	ng	96
70) Phenanthrene	11.52	178	1017277	36.792	ng	99
71) Anthracene	11.57	178	397633	14.283	ng	99
72) Carbazole	11.74	167	145477	5.025	ng	98
74) Fluoranthene	12.71	202	1085723	36.594	ng	97
77) Pyrene	12.95	202	965796	31.587	ng	97
80) Benzo(a)anthracene	14.14	228	598267	22.838	ng	93
82) Chrysene	14.18	228	594736	23.635	ng	98
85) Indeno(1,2,3-cd)pyrene	17.25	276	268544	12.592	ng	# 89
87) Benzo(b)fluoranthene	15.22	252	716362m	25.763	ng	
88) Benzo(k)fluoranthene	15.25	252	299068m	10.978	ng	
89) Benzo(a)pyrene	15.61	252	555047	21.823	ng	96
90) Dibenz(a,h)anthracene	17.25	278	62743m	2.854	ng	
91) Benzo(g,h,i)perylene	17.72	276	252792m	11.656	ng	

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

