

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071718\
 Data File : BF107449.D
 Acq On : 17 Jul 2018 15:10
 Operator : JU/SJ
 Sample : J3829-05 2X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-01-071318-C

Manual Integrations
 APPROVED

Sohil
 7/18/2018 10:44:03 AM

Quant Time: Jul 17 18:34:09 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 17 11:29:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	144731	20.00	ng	0.00
21) Naphthalene-d8	8.27	136	549073	20.00	ng	0.00
38) Acenaphthene-d10	10.03	164	273609	20.00	ng	0.00
63) Phenanthrene-d10	11.52	188	511162	20.00	ng	0.00
75) Chrysene-d12	14.18	240	362326	20.00	ng	0.00
86) Perylene-d12	15.72	264	282246	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	364165	38.21	ng	0.00
7) Phenol-d6	6.61	99	482963	39.20	ng	0.00
23) Nitrobenzene-d5	7.54	82	327393	31.80	ng	-0.01
41) 2,4,6-Tribromophenol	10.82	330	103513	43.06	ng	0.00
44) 2-Fluorobiphenyl	9.34	172	581904	28.22	ng	0.00
78) Terphenyl-d14	13.11	244	501979	34.42	ng	0.00
Target Compounds						
31) Naphthalene	8.29	128	94435	3.541	ng	97
37) 2-Methylnaphthalene	8.98	142	45331	2.579	ng	# 85
48) Acenaphthylene	9.89	152	131991	5.062	ng	95
49) Dimethylphthalate	9.74	163	111355	5.122	ng	96
51) Acenaphthene	10.06	154	75647	4.376	ng	98
54) Dibenzofuran	10.24	168	118473	4.435	ng	# 95
57) Fluorene	10.58	166	180927	10.227	ng	# 96
70) Phenanthrene	11.55	178	1188139	46.641	ng	97
71) Anthracene	11.60	178	462355	18.223	ng	97
72) Carbazole	11.75	167	96504	4.017	ng	# 86
74) Fluoranthene	12.75	202	1517382	52.407	ng	94
77) Pylene	12.98	202	1377294	55.899	ng	99
80) Benzo(a)anthracene	14.17	228	789785	35.625	ng	95
82) Chrysene	14.21	228	684296	32.313	ng	99
85) Indeno(1,2,3-cd)pyrene	17.31	276	223598	14.833	ng	# 88
87) Benzo(b)fluoranthene	15.26	252	690563m	41.801	ng	
88) Benzo(k)fluoranthene	15.29	252	283841m	17.605	ng	
89) Benzo(a)pyrene	15.65	252	497143m	32.786	ng	
90) Dibenzo(a,h)anthracene	17.31	278	58018	5.070	ng	95
91) Benzo(a,h,i)perylene	17.79	276	212340	18.261	ng	95

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

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