

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102518\
 Data File : BF110172.D
 Acq On : 25 Oct 2018 20:39
 Operator : JU/SJ
 Sample : J5625-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DW-02

Manual Integrations
 APPROVED

Sohil
 10/26/2018 3:49:22 PM

Quant Time: Oct 26 04:47:25 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 23 15:06:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	167914	20.00	ng	-0.02
21) Naphthalene-d8	8.25	136	677605	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	309279	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	507677	20.00	ng	-0.02
76) Chrysene-d12	14.14	240	302112	20.00	ng	-0.02
87) Perylene-d12	15.67	264	291672	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	1019765	98.90	ng	0.00
7) Phenol-d6	6.59	99	1295013	98.19	ng	-0.01
23) Nitrobenzene-d5	7.53	82	831103	72.75	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	271312	88.56	ng	-0.01
45) 2-Fluorobiphenyl	9.32	172	1325708	67.31	ng	-0.02
79) Terphenyl-d14	13.09	244	909060	62.58	ng	-0.02
Target Compounds						
10) Phenol	6.60	94	147051	10.431	ng	Qvalue # 62
31) Naphthalene	8.27	128	70866	2.269	ng	99
37) 2-Methylnaphthalene	8.96	142	46122	2.232	ng	100
50) Dimethylphthalate	9.72	163	256818	11.249	ng	99
71) Phenanthrene	11.52	178	200384	7.543	ng	99
72) Anthracene	11.57	178	63001	2.366	ng	99
75) Fluoranthene	12.72	202	305842	11.345	ng	100
78) Pyrene	12.95	202	346940	16.018	ng	98
80) Butylbenzylphthalate	13.55	149	90152	9.590	ng	96
81) Benzo(a)anthracene	14.13	228	176654	9.860	ng	95
83) Chrysene	14.17	228	175201	10.296	ng	98
84) Bis(2-ethylhexyl)phthalate	14.10	149	67344	5.299	ng	98
86) Indeno(1,2,3-cd)pyrene	17.23	276	100767	7.138	ng	99
88) Benzo(b)fluoranthene	15.22	252	239271m	14.206	ng	
89) Benzo(k)fluoranthene	15.24	252	90830m	5.485	ng	
90) Benzo(a)pyrene	15.60	252	158590	10.233	ng	98
91) Dibenzo(a,h)anthracene	17.23	278	29086	2.175	ng	# 77
92) Benzo(g,h,i)perylene	17.70	276	95977	7.283	ng	94

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

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