

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102518\
 Data File : BF110164.D
 Acq On : 25 Oct 2018 17:00
 Operator : JU/SJ
 Sample : J5636-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HP-GRID-A

Manual Integrations
 APPROVED

Sohil
 10/26/2018 3:49:17 PM

Quant Time: Oct 26 03:55:44 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	142045	20.00	ng	-0.02
21) Naphthalene-d8	8.25	136	564259	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	273822	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	499612	20.00	ng	-0.02
76) Chrysene-d12	14.14	240	303112	20.00	ng	-0.02
87) Perylene-d12	15.67	264	263042	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	894541	102.55	ng	0.00
7) Phenol-d6	6.59	99	1142406	102.39	ng	-0.01
23) Nitrobenzene-d5	7.53	82	716204	75.29	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	262268	96.69	ng	-0.01
45) 2-Fluorobiphenyl	9.32	172	1237242	70.95	ng	-0.02
79) Terphenyl-d14	13.09	244	1065490	73.11	ng	-0.02
Target Compounds						
10) Phenol	6.60	94	60698	5.090	ng	Qvalue # 71
50) Dimethylphthalate	9.71	163	135268	6.692	ng	99
52) Acenaphthene	10.04	154	38249	2.204	ng	98
71) Phenanthrene	11.52	178	312267	11.945	ng	98
72) Anthracene	11.57	178	371930	14.194	ng	98
75) Fluoranthene	12.72	202	234560	8.841	ng	100
78) Pyrene	12.94	202	273810	12.600	ng	99
81) Benzo(a)anthracene	14.13	228	82369m	4.582	ng	
83) Chrysene	14.17	228	90559	5.304	ng	98
86) Indeno(1,2,3-cd)pyrene	17.23	276	34014	2.402	ng	97
88) Benzo(b)fluoranthene	15.22	252	74447m	4.901	ng	
90) Benzo(a)pyrene	15.60	252	64187	4.592	ng	99
92) Benzo(a,h,i)perylene	17.70	276	35115	2.955	ng	# 95

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

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