

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102518\
 Data File : BF110168.D
 Acq On : 25 Oct 2018 18:50
 Operator : JU/SJ
 Sample : J5636-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HP-GRID-C

Manual Integrations
 APPROVED

Sohil
 10/26/2018 3:49:20 PM

Quant Time: Oct 26 04:27:20 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	174632	20.00	ng	-0.02
21) Naphthalene-d8	8.25	136	710644	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	338924	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	607542	20.00	ng	-0.02
76) Chrysene-d12	14.14	240	331879	20.00	ng	-0.02
87) Perylene-d12	15.67	264	295948	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	1298818	121.12	ng	0.00
7) Phenol-d6	6.59	99	1676400	122.22	ng	-0.01
23) Nitrobenzene-d5	7.53	82	1058008	88.31	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	372081	110.83	ng	-0.01
45) 2-Fluorobiphenyl	9.33	172	1658783	76.85	ng	-0.01
79) Terphenyl-d14	13.09	244	1361484	85.32	ng	-0.02
Target Compounds						
10) Phenol	6.60	94	97377	6.641	ng	Qvalue # 64
31) Naphthalene	8.27	128	102086	3.117	ng	99
50) Dimethylphthalate	9.71	163	173943	6.953	ng	99
52) Acenaphthene	10.04	154	71456	3.326	ng	96
58) Fluorene	10.56	166	46894	2.095	ng	97
71) Phenanthrene	11.52	178	429465	13.510	ng	99
72) Anthracene	11.57	178	138240	4.338	ng	99
75) Fluoranthene	12.72	202	301821	9.355	ng	99
78) Pyrene	12.94	202	369606	15.534	ng	99
81) Benzo(a)anthracene	14.13	228	98504m	5.005	ng	
83) Chrysene	14.17	228	86729	4.640	ng	97
86) Indeno(1,2,3-cd)pyrene	17.22	276	40120	2.587	ng	97
88) Benzo(b)fluoranthene	15.22	252	81582m	4.774	ng	
90) Benzo(a)pyrene	15.60	252	81832	5.204	ng	96
92) Benzo(a,h,i)perylene	17.70	276	41094	3.073	ng	97

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

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