

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112818\
 Data File : BF111061.D
 Acq On : 28 Nov 2018 22:29
 Operator : JU/SJ
 Sample : J6062-02DL 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-20(1-10)DL

Manual Integrations
 APPROVED

mohammad
 11/29/2018 1:58:47 PM

Quant Time: Nov 29 03:43:15 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 27 15:32:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	60432	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	239455	20.00	ng	0.00
39) Acenaphthene-d10	9.97	164	95135	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	147937	20.00	ng	0.00
76) Chrysene-d12	14.13	240	130410	20.00	ng	0.00
87) Perylene-d12	15.66	264	98010	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	163681	44.92	ng	0.00
7) Phenol-d6	6.56	99	221400	46.62	ng	0.00
23) Nitrobenzene-d5	7.50	82	130046	31.08	ng	0.00
42) 2,4,6-Tribromophenol	10.77	330	34887	36.92	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	219951	33.50	ng	0.00
79) Terphenyl-d14	13.05	244	158124	23.60	ng	-0.01
Target Compounds						
10) Phenol	6.58	94	12499	2.337	ng	93
50) Dimethylphthalate	9.69	163	15685	2.248	ng	99
55) Dibenzofuran	10.19	168	17518	2.071	ng	99
58) Fluorene	10.53	166	26213	3.984	ng	98
71) Phenanthrene	11.49	178	222459	28.392	ng	99
72) Anthracene	11.55	178	71963	9.019	ng	97
75) Fluoranthene	12.69	202	274206	34.051	ng	99
78) Pyrene	12.92	202	213017	22.364	ng	97
81) Benzo(a)anthracene	14.12	228	114984	14.934	ng	98
83) Chrysene	14.15	228	89780	11.846	ng	99
86) Indeno(1,2,3-cd)pyrene	17.26	276	29700	5.459	ng	100
88) Benzo(b)fluoranthene	15.20	252	91564m	15.903	ng	
89) Benzo(k)fluoranthene	15.23	252	27485m	4.638	ng	
90) Benzo(a)pyrene	15.60	252	57137	10.661	ng	98
92) Benzo(a,h,i)perylene	17.75	276	24964	5.855	ng	# 88

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

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