

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112718\
 Data File : BF111003.D
 Acq On : 27 Nov 2018 12:15
 Operator : JU/SJ
 Sample : J6028-06RE
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-2RE

Manual Integrations
 APPROVED

mohammad
 11/28/2018 4:42:06 PM

Quant Time: Nov 27 14:39:49 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 06:43:50 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	50888	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	184287	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	69701	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	126123	20.00	ng	0.00
76) Chrysene-d12	14.13	240	115454	20.00	ng	0.00
87) Perylene-d12	15.67	264	76468	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	317441	103.45	ng	0.01
7) Phenol-d6	6.57	99	383461	95.89	ng	0.00
23) Nitrobenzene-d5	7.50	82	243612	75.65	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	64730	93.51	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	399375	83.01	ng	0.00
79) Terphenyl-d14	13.06	244	381971	64.40	ng	0.00
Target Compounds						
10) Phenol	6.58	94	21178	4.702	ng	82
50) Dimethylphthalate	9.69	163	52357	10.242	ng	99
71) Phenanthrene	11.50	178	107414	16.080	ng	99
72) Anthracene	11.56	178	32673m	4.803	ng	
75) Fluoranthene	12.70	202	203752	29.678	ng	99
78) Pyrene	12.93	202	185219	21.965	ng	98
81) Benzo(a)anthracene	14.13	228	89009	13.058	ng	99
83) Chrysene	14.16	228	78103m	11.640	ng	
86) Indeno(1,2,3-cd)pyrene	17.27	276	29186	6.060	ng	98
88) Benzo(b)fluoranthene	15.22	252	79746m	17.752	ng	
89) Benzo(k)fluoranthene	15.24	252	23620m	5.108	ng	
90) Benzo(a)pyrene	15.61	252	50971	12.190	ng	# 96
91) Dibenzo(a,h)anthracene	17.27	278	6970	2.020	ng	# 67
92) Benzo(a,h,i)perylene	17.77	276	29747	8.942	ng	93

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

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