

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112618\
 Data File : BF110982.D
 Acq On : 26 Nov 2018 23:46
 Operator : JU/SJ
 Sample : J6028-06
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-2

Manual Integrations
 APPROVED

mohammad
 11/27/2018 11:48:16 AM

Quant Time: Nov 27 06:11:26 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 26 15:26:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	54386	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	197398	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	79341	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	140052	20.00	ng	0.00
76) Chrysene-d12	14.13	240	87369	20.00	ng	0.00
87) Perylene-d12	15.67	264	66635	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	348734	106.33	ng	0.01
7) Phenol-d6	6.57	99	426187	99.72	ng	0.00
23) Nitrobenzene-d5	7.50	82	245153	71.08	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	72225	91.66	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	440559	80.45	ng	0.00
79) Terphenyl-d14	13.06	244	362654	80.80	ng	0.00
Target Compounds						
10) Phenol	6.58	94	23905	4.967	ng	81
50) Dimethylphthalate	9.69	163	60189	10.344	ng	99
71) Phenanthrene	11.50	178	122380	16.499	ng	98
72) Anthracene	11.56	178	33266	4.404	ng	96
75) Fluoranthene	12.70	202	198450	26.031	ng	99
78) Pyrene	12.93	202	169145	26.507	ng	98
81) Benzo(a)anthracene	14.13	228	70324	13.634	ng	98
83) Chrysene	14.16	228	58793	11.579	ng	97
86) Indeno(1,2,3-cd)pyrene	17.28	276	25976	7.127	ng	94
88) Benzo(b)fluoranthene	15.22	252	63908m	16.326	ng	
89) Benzo(k)fluoranthene	15.24	252	25486m	6.325	ng	
90) Benzo(a)pyrene	15.61	252	46428	12.742	ng	# 96
91) Dibenzo(a,h)anthracene	17.27	278	6859	2.281	ng	# 73
92) Benzo(a,h,i)perylene	17.77	276	28143	9.709	ng	# 88

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

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