

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112618\
 Data File : BF110972.D
 Acq On : 26 Nov 2018 19:10
 Operator : JU/SJ
 Sample : J5996-25 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AOC1-06-A

Manual Integrations
 APPROVED

mohammad
 11/27/2018 11:47:48 AM

Quant Time: Nov 27 03:35:03 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	6.93	152	71577	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	271064	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	104304	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	168169	20.00	ng	0.00
76) Chrysene-d12	14.13	240	110666	20.00	ng	0.00
87) Perylene-d12	15.67	264	82128	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	150236	34.81	ng	0.00
7) Phenol-d6	6.56	99	285217	50.71	ng	0.00
23) Nitrobenzene-d5	7.50	82	189343	39.98	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	10198	9.84	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	324333	45.05	ng	0.00
79) Terphenyl-d14	13.06	244	256925	45.19	ng	0.00
Target Compounds						
10) Phenol	6.58	94	13730	2.167	ng	86
50) Dimethylphthalate	9.69	163	21159	2.766	ng	99
75) Fluoranthene	12.70	202	51028	5.574	ng	98
78) Pyrene	12.93	202	49214	6.089	ng	96
81) Benzo(a)anthracene	14.12	228	24141	3.695	ng	# 87
83) Chrysene	14.16	228	24990	3.885	ng	# 92
86) Indeno(1,2,3-cd)pyrene	17.29	276	12083	2.617	ng	93
88) Benzo(b)fluoranthene	15.22	252	26502m	5.493	ng	
89) Benzo(k)fluoranthene	15.24	252	10302m	2.075	ng	
90) Benzo(a)pyrene	15.61	252	19198	4.275	ng	# 73
92) Benzo(g,h,i)perylene	17.77	276	14692	4.112	ng	# 77

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

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