

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111618\  
 Data File : BF110823.D  
 Acq On : 16 Nov 2018 18:13  
 Operator : JU/SJ  
 Sample : J5952-06 5X  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampled :  
 DUP-01

Manual Integrations  
 APPROVED

Sohil  
 11/19/2018 11:55:58 AM

Quant Time: Nov 16 23:46:05 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Nov 13 12:43:10 2018  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	61603	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	257987	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	116669	20.00	ng	-0.01
64) Phenanthrene-d10	11.47	188	184799	20.00	ng	-0.01
76) Chrysene-d12	14.13	240	146450	20.00	ng	-0.01
87) Perylene-d12	15.66	264	112745	20.00	ng	0.00

## System Monitoring Compounds

5) 2-Fluorophenol	5.56	112	98438	25.96	ng	0.00
7) Phenol-d6	6.56	99	123797	25.53	ng	-0.01
23) Nitrobenzene-d5	7.50	82	75093	16.76	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	24105	20.50	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	141693	17.35	ng	-0.01
79) Terphenyl-d14	13.06	244	101338	12.96	ng	-0.01

## Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
31) Naphthalene	8.25	128	45860	3.787	ng	99
49) Acenaphthylene	9.85	152	60190	5.330	ng	99
50) Dimethylphthalate	9.69	163	19545	2.245	ng	99
55) Dibenzofuran	10.19	168	25423	2.435	ng	95
58) Fluorene	10.53	166	45363	5.657	ng	95
71) Phenanthrene	11.50	178	324318	32.757	ng	100
72) Anthracene	11.55	178	119271	11.942	ng	98
75) Fluoranthene	12.70	202	401701	40.470	ng	98
78) Pyrene	12.93	202	396584	35.170	ng	97
81) Benzo(a)anthracene	14.12	228	232127	26.518	ng	96
83) Chrysene	14.16	228	216460	25.956	ng	97
86) Indeno(1,2,3-cd)pyrene	17.23	276	70549	11.789	ng	99
88) Benzo(b)fluoranthene	15.20	252	239925m	35.572	ng	
89) Benzo(k)fluoranthene	15.23	252	74843m	11.230	ng	
90) Benzo(a)pyrene	15.59	252	160205	26.143	ng	99
91) Dibenzo(a,h)anthracene	17.23	278	21273m	4.102	ng	
92) Benzo(a,h,i)perylene	17.71	276	71939	13.981	ng	98

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

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