

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071318\
 Data File : BF107371.D
 Acq On : 13 Jul 2018 17:02
 Operator : JU/SJ
 Sample : J3825-09 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-071118-E

Manual Integrations
 APPROVED

Sohil
 7/16/2018 3:22:31 PM

Quant Time: Jul 13 19:05:41 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 29 14:38:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	53757	20.00	ng	-0.03
21) Naphthalene-d8	8.21	136	192306	20.00	ng	-0.04
38) Acenaphthene-d10	9.98	164	89191	20.00	ng	-0.03
63) Phenanthrene-d10	11.48	188	177673	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	152940	20.00	ng	-0.03
86) Perylene-d12	15.68	264	115769	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	169468	53.38	ng	-0.03
7) Phenol-d6	6.58	99	203120	51.23	ng	-0.04
23) Nitrobenzene-d5	7.50	82	118704	34.30	ng	-0.04
41) 2,4,6-Tribromophenol	10.78	330	51827	50.14	ng	-0.02
44) 2-Fluorobiphenyl	9.29	172	231789	35.38	ng	-0.03
78) Terphenyl-d14	13.06	244	281101	44.52	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.68	163	31857	4.762	ng	Qvalue # 97
57) Fluorene	10.53	166	19184	3.174	ng	99
70) Phenanthrene	11.50	178	143046	16.692	ng	98
71) Anthracene	11.55	178	47408	5.420	ng	98
74) Fluoranthene	12.70	202	177511	18.793	ng	93
77) Pyrene	12.93	202	152505	15.121	ng	99
80) Benzo(a)anthracene	14.12	228	68203	7.317	ng	97
82) Chrysene	14.16	228	58849	6.862	ng	96
85) Indeno(1,2,3-cd)pyrene	17.28	276	24194	3.325	ng	95
87) Benzo(b)fluoranthene	15.22	252	60948m	8.475	ng	
88) Benzo(k)fluoranthene	15.25	252	17893m	2.613	ng	
89) Benzo(a)pyrene	15.62	252	41022	6.417	ng	95
91) Benzo(a,h,i)perylene	17.77	276	25603	4.835	ng	# 87

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

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