

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080318\
 Data File : BF107971.D
 Acq On : 4 Aug 2018 6:48
 Operator : JU/SJ
 Sample : J4262-07DL 2X
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUP-073118DL

Manual Integrations
 APPROVED

Sohil
 8/6/2018 5:06:16 PM

Quant Time: Aug 06 04:16:27 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 30 17:48:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	65324	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	296446	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	120351	20.00	ng	-0.01
63) Phenanthrene-d10	11.49	188	246001	20.00	ng	0.00
75) Chrysene-d12	14.15	240	247985	20.00	ng	0.00
86) Perylene-d12	15.68	264	179427	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	39401	10.25	ng	0.04
7) Phenol-d6	6.64	99	51189	10.11	ng	0.03
23) Nitrobenzene-d5	7.52	82	171081	33.25	ng	-0.01
41) 2,4,6-Tribromophenol	10.79	330	89985	55.03	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	422056	48.87	ng	-0.01
78) Terphenyl-d14	13.08	244	480786	42.04	ng	0.00
Target Compounds						
27) 2,4-Dimethylphenol	7.92	122	7389	2.038	ng	Qvalue # 82
31) Naphthalene	8.26	128	811289	58.874	ng	100
37) 2-Methylnaphthalene	8.95	142	70309	8.123	ng	94
48) Acenaphthylene	9.86	152	34094	2.663	ng	# 92
51) Acenaphthene	10.02	154	129578	17.421	ng	100
54) Dibenzofuran	10.21	168	79618	6.925	ng	98
57) Fluorene	10.55	166	103646	11.600	ng	98
70) Phenanthrene	11.52	178	107546	9.207	ng	98
71) Anthracene	11.58	178	31517m	2.339	ng	
72) Carbazole	11.74	167	195781	14.704	ng	97
74) Fluoranthene	12.71	202	29336	2.138	ng	92

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

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