

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072220\
 Data File : BM026842.D
 Acq On : 22 Jul 2020 21:07
 Operator : JU/CG
 Sample : L3360-04RX
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SP-3RX

Manual Integrations
 APPROVED

Sohil
 7/24/2020 12:11:43 PM

Quant Time: Jul 23 03:09:25 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM070820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 22 14:16:10 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.30	152	97230	20.00	ng	0.00
21) Naphthalene-d8	10.05	136	332901	20.00	ng	0.00
39) Acenaphthene-d10	13.96	164	183643	20.00	ng	0.00
64) Phenanthrene-d10	16.72	188	389290	20.00	ng	0.00
76) Chrysene-d12	20.95	240	444654	20.00	ng	0.00
86) Perylene-d12	23.01	264	464381	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.94	112	82646	14.25	ng	0.00
7) Phenol-d6	6.51	99	384365	41.59	ng	0.00
23) Nitrobenzene-d5	8.45	82	400898	51.38	ng	0.00
42) 2,4,6-Tribromophenol	15.47	330	20356	7.61	ng	0.00
45) 2-Fluorobiphenyl	12.57	172	672664	48.70	ng	0.00
79) Terphenyl-d14	19.39	244	992185	43.91	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.44	163	73143	4.657	ng	Qvalue 100
71) Phenanthrene	16.77	178	372876	16.225	ng	99
72) Anthracene	16.85	178	89776	3.924	ng	100
73) Carbazole	17.14	167	52280	2.397	ng	98
75) Fluoranthene	18.79	202	1042448	36.828	ng	98
78) Pyrene	19.16	202	936273	33.425	ng	99
81) Benzo(a)anthracene	20.93	228	545731	18.271	ng	98
83) Chrysene	20.98	228	521633	17.943	ng	97
87) Indeno(1,2,3-cd)pyrene	25.01	276	289807	8.766	ng	95
88) Benzo(b)fluoranthene	22.41	252	625241m	20.434	ng	
89) Benzo(k)fluoranthene	22.45	252	229648m	7.652	ng	
90) Benzo(a)pyrene	22.92	252	453444	15.963	ng	# 95
91) Dibenzo(a,h)anthracene	25.01	278	76125	2.728	ng	# 76
92) Benzo(g,h,i)perylene	25.61	276	287768	11.191	ng	98

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

